

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081325\
 Data File : BP025389.D
 Acq On : 13 Aug 2025 18:13
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 08/14/2025
 Supervised By :Jagrut Upadhyay 08/14/2025

Quant Time: Aug 13 18:54:49 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 06 06:41:44 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.801	152	184702	20.000	ng	0.00	
21) Naphthalene-d8	10.578	136	736218	20.000	ng	0.00	
39) Acenaphthene-d10	14.424	164	463395	20.000	ng	0.00	
64) Phenanthrene-d10	17.230	188	942404	20.000	ng	0.00	
76) Chrysene-d12	21.665	240	1019671	20.000	ng	0.00	
86) Perylene-d12	25.059	264	1166515	20.000	ng	0.02	
System Monitoring Compounds							
5) 2-Fluorophenol	5.413	112	896126	78.034	ng	0.00	
7) Phenol-d6	6.972	99	1100976	72.847	ng	0.00	
23) Nitrobenzene-d5	8.936	82	1136884	85.552	ng	0.00	
42) 2,4,6-Tribromophenol	15.936	330	512553	110.873	ng	0.01	
45) 2-Fluorobiphenyl	13.036	172	2714182	80.900	ng	0.00	
79) Terphenyl-d14	19.948	244	4405762	81.610	ng	0.00	
Target Compounds							
2) 1,4-Dioxane	3.331	88	171067	37.748	ng	96	Qvalue
3) Pyridine	3.731	79	442368	35.134	ng	96	
4) n-Nitrosodimethylamine	3.631	42	188153	32.188	ng	95	#
6) Aniline	7.131	93	724014	34.897	ng	99	
8) 2-Chlorophenol	7.372	128	500853	39.844	ng	97	
9) Benzaldehyde	6.942	77	448384	41.771	ng	98	
10) Phenol	7.001	94	568907	35.181	ng	98	
11) bis(2-Chloroethyl)ether	7.219	93	457720	36.125	ng	98	
12) 1,3-Dichlorobenzene	7.689	146	561340	39.743	ng	99	
13) 1,4-Dichlorobenzene	7.831	146	570722	39.939	ng	97	
14) 1,2-Dichlorobenzene	8.148	146	543452	39.259	ng	98	
15) Benzyl Alcohol	8.031	79	412992	35.252	ng	99	
16) 2,2'-oxybis(1-Chloropr...	8.313	45	628261	31.389	ng	95	
17) 2-Methylphenol	8.231	107	392344	36.260	ng	97	
18) Hexachloroethane	8.872	117	213457	41.770	ng	91	
19) n-Nitroso-di-n-propyla...	8.589	70	346148	34.859	ng	94	
20) 3+4-Methylphenols	8.554	107	525616	35.276	ng	99	
22) Acetophenone	8.601	105	716512	38.768	ng	97	
24) Nitrobenzene	8.978	77	507543	40.347	ng	96	
25) Isophorone	9.513	82	982573	37.145	ng	99	
26) 2-Nitrophenol	9.684	139	271753	49.472	ng	97	
27) 2,4-Dimethylphenol	9.748	122	455846	37.378	ng	94	
28) bis(2-Chloroethoxy)met...	9.978	93	613737	38.140	ng	100	
29) 2,4-Dichlorophenol	10.225	162	450740	42.216	ng	99	
30) 1,2,4-Trichlorobenzene	10.436	180	508300	43.152	ng	99	
31) Naphthalene	10.625	128	1527522	39.968	ng	99	
32) Benzoic acid	9.883	122	318796	46.453	ng	99	
33) 4-Chloroaniline	10.725	127	625936	36.824	ng	99	
34) Hexachlorobutadiene	10.913	225	322850	47.747	ng	98	
35) Caprolactam	11.507	113	153606	37.085	ng	87	#
36) 4-Chloro-3-methylphenol	11.854	107	486095	39.346	ng	99	
37) 2-Methylnaphthalene	12.230	142	985811	39.851	ng	100	
38) 1-Methylnaphthalene	12.454	142	1027342	39.802	ng	100	
40) 1,2,4,5-Tetrachloroben...	12.607	216	559369	44.214	ng	99	
41) Hexachlorocyclopentadiene	12.595	237	299470	38.353	ng	98	
43) 2,4,6-Trichlorophenol	12.848	196	368871	45.134	ng	99	

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081325\
 Data File : BP025389.D
 Acq On : 13 Aug 2025 18:13
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
ClientSampleId :
 SSTDCCC040EC

Manual Integrations
APPROVED
 Reviewed By :Rahul Chavli 08/14/2025
 Supervised By :Jagrut Upadhyay 08/14/2025

Quant Time: Aug 13 18:54:49 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 06 06:41:44 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.919	196	402704	43.925	ng	99
46) 1,1'-Biphenyl	13.254	154	1363544	39.928	ng	99
47) 2-Chloronaphthalene	13.295	162	1065180	40.884	ng	99
48) 2-Nitroaniline	13.483	65	300035	39.849	ng	98
49) Acenaphthylene	14.136	152	1721913	39.622	ng	100
50) Dimethylphthalate	13.872	163	1341204	39.703	ng	100
51) 2,6-Dinitrotoluene	13.989	165	287495	46.139	ng	92
52) Acenaphthene	14.483	154	1084324m	40.905	ng	
53) 3-Nitroaniline	14.319	138	309614	42.464	ng	94
54) 2,4-Dinitrophenol	14.524	184	117674	46.258	ng #	50
55) Dibenzofuran	14.824	168	1594419	40.249	ng	99
56) 4-Nitrophenol	14.636	139	253454	39.388	ng	96
57) 2,4-Dinitrotoluene	14.783	165	413517	43.163	ng	99
58) Fluorene	15.489	166	1236568	39.744	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.060	232	352030	45.967	ng	97
60) Diethylphthalate	15.260	149	1405313	41.263	ng	100
61) 4-Chlorophenyl-phenyle...	15.483	204	631982	42.846	ng	99
62) 4-Nitroaniline	15.495	138	309199	42.799	ng	98
63) Azobenzene	15.783	77	1193034	37.708	ng	98
65) 4,6-Dinitro-2-methylph...	15.560	198	187566	45.770	ng	94
66) n-Nitrosodiphenylamine	15.701	169	1127513	39.294	ng	100
67) 4-Bromophenyl-phenylether	16.401	248	412364	43.488	ng	98
68) Hexachlorobenzene	16.518	284	474025	44.553	ng	95
69) Atrazine	16.677	200	428489	42.462	ng	99
70) Pentachlorophenol	16.871	266	316137	46.638	ng	99
71) Phenanthrene	17.271	178	2027625	39.282	ng	100
72) Anthracene	17.365	178	2100517	40.360	ng	100
73) Carbazole	17.636	167	1967991	39.392	ng	99
74) Di-n-butylphthalate	18.248	149	2689910	44.506	ng	99
75) Fluoranthene	19.348	202	2507136	41.959	ng	98
77) Benzidine	19.542	184	1560704	42.288	ng	99
78) Pyrene	19.724	202	2631764	38.813	ng	99
80) Butylbenzylphthalate	20.677	149	1300248	47.282	ng	96
81) Benzo(a)anthracene	21.648	228	2703813	39.903	ng	100
82) 3,3'-Dichlorobenzidine	21.571	252	1031182	43.133	ng	98
83) Chrysene	21.712	228	2508190	39.909	ng	99
84) Bis(2-ethylhexyl)phtha...	21.601	149	2030394	46.239	ng	99
85) Di-n-octyl phthalate	22.900	149	3513733	47.807	ng	96
87) Indeno(1,2,3-cd)pyrene	28.947	276	3398258	40.644	ng	99
88) Benzo(b)fluoranthene	23.995	252	2746202	39.327	ng	99
89) Benzo(k)fluoranthene	24.071	252	2706136	39.093	ng	98
90) Benzo(a)pyrene	24.906	252	2669640	39.699	ng	99
91) Dibenzo(a,h)anthracene	29.024	278	2785665	40.680	ng	97
92) Benzo(g,h,i)perylene	30.124	276	2693148	40.074	ng	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081325\
 Data File : BP025389.D
 Acq On : 13 Aug 2025 18:13
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
ClientSampleId :
 SSTDCCC040EC

Quant Time: Aug 13 18:54:49 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 06 06:41:44 2025
 Response via : Initial Calibration

Manual Integrations
APPROVED
 Reviewed By :Rahul Chavli 08/14/2025
 Supervised By :Jagrut Upadhyay 08/14/2025

