

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080725\
 Data File : BP025313.D
 Acq On : 07 Aug 2025 14:28
 Operator : CG/JU
 Sample : Q2763-04MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-2MS

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 08/08/2025
 Supervised By :Jagrut Upadhyay 08/11/2025

Quant Time: Aug 07 14:46:07 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	7.802	152	221339	20.000	ng	0.00	
21) Naphthalene-d8	10.572	136	916218	20.000	ng	0.00	
39) Acenaphthene-d10	14.419	164	588971	20.000	ng	0.00	
64) Phenanthrene-d10	17.213	188	1201001	20.000	ng	-0.01	
76) Chrysene-d12	21.660	240	1176296	20.000	ng	0.00	
86) Perylene-d12	25.036	264	1330316	20.000	ng	0.00	
System Monitoring Compounds							
5) 2-Fluorophenol	5.414	112	1562859	113.565	ng	0.00	
7) Phenol-d6	6.972	99	1900658	104.942	ng	0.00	
23) Nitrobenzene-d5	8.943	82	1314181	79.465	ng	0.00	
42) 2,4,6-Tribromophenol	15.931	330	834058	141.951	ng	0.00	
45) 2-Fluorobiphenyl	13.031	172	3029872	71.054	ng	0.00	
79) Terphenyl-d14	19.942	244	4718093	75.759	ng	0.00	
Target Compounds							
2) 1,4-Dioxane	3.349	88	181515	33.424	ng		Qvalue # 68
3) Pyridine	3.737	79	496876	32.931	ng		97
4) n-Nitrosodimethylamine	3.643	42	249221	35.578	ng		95
6) Aniline	7.137	93	544442	21.898	ng		97
8) 2-Chlorophenol	7.372	128	639740	42.469	ng		99
9) Benzaldehyde	6.949	77	281865	21.912	ng		96
10) Phenol	6.996	94	721038	37.208	ng		100
11) bis(2-Chloroethyl)ether	7.231	93	572421	37.700	ng		99
12) 1,3-Dichlorobenzene	7.696	146	637576	37.668	ng		99
13) 1,4-Dichlorobenzene	7.837	146	648981	37.899	ng		98
14) 1,2-Dichlorobenzene	8.155	146	627169	37.808	ng		99
15) Benzyl Alcohol	8.031	79	567057	40.391	ng		99
16) 2,2'-oxybis(1-Chloropr...	8.331	45	835634	34.839	ng		99
17) 2-Methylphenol	8.231	107	537504	41.453	ng		99
18) Hexachloroethane	8.878	117	231212	37.755	ng		98
19) n-Nitroso-di-n-propyla...	8.596	70	449798	37.799	ng		98
20) 3+4-Methylphenols	8.549	107	727470	40.742	ng		97
22) Acetophenone	8.608	105	918172	39.920	ng		99
24) Nitrobenzene	8.984	77	648775	41.442	ng		95
25) Isophorone	9.508	82	1319191	40.073	ng		100
26) 2-Nitrophenol	9.684	139	319115	46.886	ng		98
27) 2,4-Dimethylphenol	9.749	122	641648	42.277	ng		98
28) bis(2-Chloroethoxy)met...	9.984	93	809246	40.410	ng		99
29) 2,4-Dichlorophenol	10.219	162	626440	47.145	ng		99
30) 1,2,4-Trichlorobenzene	10.437	180	588995	40.179	ng		99
31) Naphthalene	10.625	128	1900022	39.947	ng		100
32) Benzoic acid	9.866	122	454934	51.160	ng		98
33) 4-Chloroaniline	10.719	127	211234	9.986	ng		99
34) Hexachlorobutadiene	10.919	225	336435	39.981	ng		97
35) Caprolactam	11.484	113	209611	40.664	ng		93
36) 4-Chloro-3-methylphenol	11.837	107	718893	46.757	ng		100
37) 2-Methylnaphthalene	12.231	142	1269364	41.232	ng		100
38) 1-Methylnaphthalene	12.449	142	1319234	41.069	ng		100
40) 1,2,4,5-Tetrachloroben...	12.602	216	668645	41.583	ng		100
41) Hexachlorocyclopentadiene	12.590	237	499288	50.310	ng		98
43) 2,4,6-Trichlorophenol	12.831	196	509284	49.029	ng		97

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080725\
 Data File : BP025313.D
 Acq On : 07 Aug 2025 14:28
 Operator : CG/JU
 Sample : Q2763-04MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
ClientSampleId :
 TP-2MS

Manual Integrations
APPROVED
 Reviewed By :Rahul Chavli 08/08/2025
 Supervised By :Jagrut Upadhyay 08/11/2025

Quant Time: Aug 07 14:46:07 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	12.907	196	581255	49.882	ng	100
46) 1,1'-Biphenyl	13.249	154	1829065	42.140	ng	99
47) 2-Chloronaphthalene	13.284	162	1388405	41.928	ng	98
48) 2-Nitroaniline	13.478	65	461747	47.626	ng	94
49) Acenaphthylene	14.131	152	2376665	43.028	ng	99
50) Dimethylphthalate	13.872	163	1922717	44.782	ng	100
51) 2,6-Dinitrotoluene	13.984	165	405340	51.182	ng	92
52) Acenaphthene	14.478	154	1478244	43.875	ng	99
53) 3-Nitroaniline	14.307	138	251805	27.172	ng	99
54) 2,4-Dinitrophenol	14.507	184	151365	46.627	ng #	59
55) Dibenzofuran	14.819	168	2165553	43.011	ng	100
56) 4-Nitrophenol	14.619	139	766453	93.714	ng	97
57) 2,4-Dinitrotoluene	14.772	165	573519	46.803	ng	99
58) Fluorene	15.478	166	1725371	43.631	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.048	232	494062	50.758	ng	98
60) Diethylphthalate	15.260	149	1989777	45.968	ng	99
61) 4-Chlorophenyl-phenyle...	15.484	204	814050	43.423	ng	98
62) 4-Nitroaniline	15.490	138	421374	45.890	ng	97
63) Azobenzene	15.778	77	1753016	43.594	ng	99
65) 4,6-Dinitro-2-methylph...	15.548	198	146153	32.045	ng	99
66) n-Nitrosodiphenylamine	15.695	169	1612540	44.097	ng	99
67) 4-Bromophenyl-phenylether	16.395	248	538797	44.587	ng	96
68) Hexachlorobenzene	16.507	284	600933	44.319	ng	98
69) Atrazine	16.678	200	582911	45.327	ng	99
70) Pentachlorophenol	16.854	266	857117	99.219	ng	100
71) Phenanthrene	17.260	178	2887546	43.897	ng	99
72) Anthracene	17.360	178	2959175	44.616	ng	100
73) Carbazole	17.625	167	2824703	44.366	ng	99
74) Di-n-butylphthalate	18.237	149	3542540	45.993	ng	99
75) Fluoranthene	19.336	202	3339110	43.850	ng	99
77) Benzidine	19.536	184	1085842	25.504	ng	99
78) Pyrene	19.713	202	3424311	43.777	ng	99
80) Butylbenzylphthalate	20.683	149	1623760	51.184	ng	99
81) Benzo(a)anthracene	21.642	228	3395999	43.445	ng	98
82) 3,3'-Dichlorobenzidine	21.566	252	975536	35.372	ng	99
83) Chrysene	21.713	228	3183194	43.906	ng	100
84) Bis(2-ethylhexyl)phtha...	21.607	149	2409679	47.570	ng	99
85) Di-n-octyl phthalate	22.913	149	4124691	48.647	ng	98
87) Indeno(1,2,3-cd)pyrene	28.877	276	4138140m	43.399	ng	
88) Benzo(b)fluoranthene	23.983	252	3378818	42.429	ng	100
89) Benzo(k)fluoranthene	24.048	252	3381495	42.834	ng	99
90) Benzo(a)pyrene	24.883	252	3316495	43.245	ng	99
91) Dibenzo(a,h)anthracene	28.995	278	3417760	43.766	ng	97
92) Benzo(g,h,i)perylene	30.089	276	3304761	43.119	ng	99

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080725\
 Data File : BP025313.D
 Acq On : 07 Aug 2025 14:28
 Operator : CG/JU
 Sample : Q2763-04MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
ClientSampleId :
 TP-2MS

Quant Time: Aug 07 14:46:07 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 08/08/2025
 Supervised By :Jagrut Upadhyay 08/11/2025

