

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101525\
 Data File : BP025979.D
 Acq On : 15 Oct 2025 18:35
 Operator : RC/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC080

Quant Time: Oct 15 23:12:33 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP101525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 15 23:04:36 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.696	152	189087	20.000	ng	0.00
21) Naphthalene-d8	10.460	136	853829	20.000	ng	0.00
39) Acenaphthene-d10	14.319	164	573100	20.000	ng	0.00
64) Phenanthrene-d10	17.125	188	1099586	20.000	ng	0.00
76) Chrysene-d12	21.577	240	809390	20.000	ng	0.00
86) Perylene-d12	24.918	264	924612	20.000	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.337	112	1821228	157.914	ng	0.00
7) Phenol-d6	6.908	99	2602237	161.683	ng	0.00
23) Nitrobenzene-d5	8.855	82	2906298	152.995	ng	0.00
42) 2,4,6-Tribromophenol	15.848	330	1247173	161.650	ng	0.00
45) 2-Fluorobiphenyl	12.931	172	6435200	143.923	ng	0.00
79) Terphenyl-d14	19.854	244	7917555	156.336	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	3.249	88	319965	70.040 ng	99
3) Pyridine	3.655	79	1002619	78.649 ng	99
4) n-Nitrosodimethylamine	3.561	42	434471	77.768 ng	100
6) Aniline	7.043	93	1750519	81.242 ng	99
8) 2-Chlorophenol	7.278	128	1067041	79.835 ng	99
9) Benzaldehyde	6.861	77	574464	66.887 ng	98
10) Phenol	6.937	94	1362857	83.868 ng	98
11) bis(2-Chloroethyl)ether	7.137	93	1051533	79.361 ng	99
12) 1,3-Dichlorobenzene	7.590	146	1118984	75.514 ng	99
13) 1,4-Dichlorobenzene	7.737	146	1139668	75.487 ng	100
14) 1,2-Dichlorobenzene	8.043	146	1126448	76.269 ng	100
15) Benzyl Alcohol	7.949	79	1106898	80.802 ng	99
16) 2,2'-oxybis(1-Chloropr...	8.219	45	1562552	76.581 ng	99
17) 2-Methylphenol	8.160	107	956854	82.151 ng	99
18) Hexachloroethane	8.760	117	435195	75.537 ng	99
19) n-Nitroso-di-n-propyla...	8.508	70	908299	79.165 ng	98
20) 3+4-Methylphenols	8.490	107	1290002	79.487 ng	97
22) Acetophenone	8.519	105	1761662	75.893 ng	# 99
24) Nitrobenzene	8.896	77	1311475	77.636 ng	98
25) Isophorone	9.425	82	2634496	77.219 ng	100
26) 2-Nitrophenol	9.596	139	668745	81.745 ng	98
27) 2,4-Dimethylphenol	9.666	122	1102040	79.495 ng	99
28) bis(2-Chloroethoxy)met...	9.884	93	1539493	77.137 ng	98
29) 2,4-Dichlorophenol	10.143	162	1132443	81.983 ng	98
30) 1,2,4-Trichlorobenzene	10.331	180	1181311	75.540 ng	97
31) Naphthalene	10.513	128	3444501	74.784 ng	100
32) Benzoic acid	9.943	122	861099	85.553 ng	96
33) 4-Chloroaniline	10.631	127	1569828	80.257 ng	98
34) Hexachlorobutadiene	10.790	225	779931	75.359 ng	100
35) Caprolactam	11.490	113	389337	79.204 ng	95
36) 4-Chloro-3-methylphenol	11.784	107	1299013	80.101 ng	98
37) 2-Methylnaphthalene	12.125	142	2280979	75.638 ng	99
38) 1-Methylnaphthalene	12.343	142	2398064	75.390 ng	100
40) 1,2,4,5-Tetrachloroben...	12.496	216	1431042	76.683 ng	100
41) Hexachlorocyclopentadiene	12.466	237	879241	79.939 ng	100
43) 2,4,6-Trichlorophenol	12.754	196	962749	81.277 ng	98

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101525\
 Data File : BP025979.D
 Acq On : 15 Oct 2025 18:35
 Operator : RC/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC080

Quant Time: Oct 15 23:12:33 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP101525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 15 23:04:36 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.848	196	1052298	81.582	ng	99
46) 1,1'-Biphenyl	13.143	154	3239748	74.836	ng	99
47) 2-Chloronaphthalene	13.190	162	2557460	75.554	ng	100
48) 2-Nitroaniline	13.407	65	868144	82.291	ng	99
49) Acenaphthylene	14.037	152	4135556	75.562	ng	100
50) Dimethylphthalate	13.784	163	3291729	75.990	ng	100
51) 2,6-Dinitrotoluene	13.901	165	733002	78.371	ng	100
52) Acenaphthene	14.384	154	2361915	74.733	ng	100
53) 3-Nitroaniline	14.248	138	766639	83.284	ng	99
54) 2,4-Dinitrophenol	14.472	184	496006	78.171	ng	95
55) Dibenzofuran	14.731	168	3754843	74.197	ng	99
56) 4-Nitrophenol	14.607	139	516599	91.443	ng	98
57) 2,4-Dinitrotoluene	14.713	165	1000653	78.959	ng	97
58) Fluorene	15.389	166	3035330	74.644	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.972	232	929263	80.289	ng	100
60) Diethylphthalate	15.160	149	3237994	75.812	ng	99
61) 4-Chlorophenyl-phenyle...	15.378	204	1612838	75.425	ng	100
62) 4-Nitroaniline	15.442	138	698625	82.358	ng	98
63) Azobenzene	15.678	77	3039918	76.360	ng	98
65) 4,6-Dinitro-2-methylph...	15.501	198	660428	83.102	ng	97
66) n-Nitrosodiphenylamine	15.601	169	2665421	75.342	ng	100
67) 4-Bromophenyl-phenylether	16.289	248	1064527	78.435	ng	96
68) Hexachlorobenzene	16.419	284	1185605	77.869	ng	98
69) Atrazine	16.584	200	981431	76.899	ng	98
70) Pentachlorophenol	16.784	266	683260	84.702	ng	98
71) Phenanthrene	17.172	178	4551842	73.039	ng	100
72) Anthracene	17.266	178	4682643	73.557	ng	100
73) Carbazole	17.554	167	4116696	73.357	ng	99
74) Di-n-butylphthalate	18.131	149	5042014	73.006	ng	100
75) Fluoranthene	19.266	202	4842528	68.886	ng	98
77) Benzidine	19.460	184	1982980	76.462	ng	100
78) Pyrene	19.642	202	4796481	80.348	ng	100
80) Butylbenzylphthalate	20.577	149	1860728	78.326	ng	98
81) Benzo(a)anthracene	21.554	228	4208838	76.030	ng	99
82) 3,3'-Dichlorobenzidine	21.477	252	1494723	75.874	ng	99
83) Chrysene	21.624	228	3961327	74.840	ng	99
84) Bis(2-ethylhexyl)phtha...	21.477	149	2498012	76.168	ng	99
85) Di-n-octyl phthalate	22.742	149	4469657	79.816	ng	99
87) Indeno(1,2,3-cd)pyrene	28.724	276	5492824	78.864	ng	# 88
88) Benzo(b)fluoranthene	23.854	252	4325350	75.113	ng	100
89) Benzo(k)fluoranthene	23.924	252	4286887	73.354	ng	100
90) Benzo(a)pyrene	24.765	252	4247050	75.907	ng	100
91) Dibenzo(a,h)anthracene	28.783	278	4509812	79.430	ng	99
92) Benzo(g,h,i)perylene	29.912	276	4440642	79.824	ng	98

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101525\
 Data File : BP025979.D
 Acq On : 15 Oct 2025 18:35
 Operator : RC/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC080

Quant Time: Oct 15 23:12:33 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP101525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 15 23:04:36 2025
 Response via : Initial Calibration

