

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060625\
 Data File : BP024867.D
 Acq On : 06 Jun 2025 15:18
 Operator : RC/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC080

Quant Time: Jun 06 16:05:27 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060625.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 06 15:53:04 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.614	152	251656	20.000	ng	0.00
21) Naphthalene-d8	10.384	136	1097725	20.000	ng	0.00
39) Acenaphthene-d10	14.254	164	655445	20.000	ng	0.00
64) Phenanthrene-d10	17.054	188	1282325	20.000	ng	0.00
76) Chrysene-d12	21.489	240	1346775	20.000	ng	0.00
86) Perylene-d12	24.742	264	1568664	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.243	112	2447087	162.312	ng	0.00
7) Phenol-d6	6.819	99	3149578	157.892	ng	0.00
23) Nitrobenzene-d5	8.766	82	3363733	148.903	ng	0.00
42) 2,4,6-Tribromophenol	15.778	330	1494447	164.916	ng	0.00
45) 2-Fluorobiphenyl	12.866	172	7387084	151.825	ng	0.00
79) Terphenyl-d14	19.789	244	11194678	148.975	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.173	88	520741	78.498	ng	100
3) Pyridine	3.567	79	1323253	82.944	ng	99
4) n-Nitrosodimethylamine	3.484	42	528461	81.024	ng	100
6) Aniline	6.961	93	2044864	80.328	ng	100
8) 2-Chlorophenol	7.196	128	1357118	79.405	ng	99
9) Benzaldehyde	6.772	77	650456	56.412	ng	99
10) Phenol	6.843	94	1640201	79.761	ng	99
11) bis(2-Chloroethyl)ether	7.049	93	1254500	77.563	ng	99
12) 1,3-Dichlorobenzene	7.502	146	1480514	77.650	ng	99
13) 1,4-Dichlorobenzene	7.649	146	1497910	77.862	ng	98
14) 1,2-Dichlorobenzene	7.961	146	1431405	75.771	ng	98
15) Benzyl Alcohol	7.866	79	1218925	79.620	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.143	45	1584064	74.829	ng	98
17) 2-Methylphenol	8.072	107	1128854	78.611	ng	99
18) Hexachloroethane	8.672	117	566168	78.169	ng	97
19) n-Nitroso-di-n-propyla...	8.419	70	1035677	76.374	ng	99
20) 3+4-Methylphenols	8.402	107	1524653	77.821	ng	99
22) Acetophenone	8.431	105	2033065	73.305	ng	# 99
24) Nitrobenzene	8.808	77	1489831	74.236	ng	99
25) Isophorone	9.331	82	3090456	78.936	ng	98
26) 2-Nitrophenol	9.508	139	870704	87.932	ng	98
27) 2,4-Dimethylphenol	9.584	122	1396457	82.403	ng	98
28) bis(2-Chloroethoxy)met...	9.808	93	1828377	78.289	ng	98
29) 2,4-Dichlorophenol	10.055	162	1419757	87.313	ng	98
30) 1,2,4-Trichlorobenzene	10.243	180	1542041	84.080	ng	99
31) Naphthalene	10.431	128	4104122	72.957	ng	100
32) Benzoic acid	9.802	122	1065621	94.314	ng	97
33) 4-Chloroaniline	10.555	127	1816811	77.087	ng	100
34) Hexachlorobutadiene	10.713	225	830962	75.174	ng	99
35) Caprolactam	11.366	113	457044	76.358	ng	97
36) 4-Chloro-3-methylphenol	11.702	107	1442747	76.925	ng	99
37) 2-Methylnaphthalene	12.049	142	2642800	74.062	ng	99
38) 1-Methylnaphthalene	12.272	142	2821350	73.955	ng	99
40) 1,2,4,5-Tetrachloroben...	12.425	216	1472269	79.160	ng	100
41) Hexachlorocyclopentadiene	12.396	237	1031961	90.970	ng	99
43) 2,4,6-Trichlorophenol	12.678	196	1039009	83.186	ng	98

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060625\
 Data File : BP024867.D
 Acq On : 06 Jun 2025 15:18
 Operator : RC/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC080

Quant Time: Jun 06 16:05:27 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060625.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 06 15:53:04 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.760	196	1117015	83.507	ng	99
46) 1,1'-Biphenyl	13.072	154	3678354	77.223	ng	99
47) 2-Chloronaphthalene	13.119	162	2834595	77.428	ng	99
48) 2-Nitroaniline	13.337	65	930628	82.685	ng	96
49) Acenaphthylene	13.966	152	4731215	77.508	ng	100
50) Dimethylphthalate	13.719	163	3769405	77.959	ng	100
51) 2,6-Dinitrotoluene	13.837	165	829970	79.625	ng	95
52) Acenaphthene	14.319	154	2715693	77.638	ng	100
53) 3-Nitroaniline	14.178	138	914372	84.530	ng	98
54) 2,4-Dinitrophenol	14.390	184	537870	79.298	ng	99
55) Dibenzofuran	14.654	168	4234879	75.424	ng	99
56) 4-Nitrophenol	14.531	139	720226	78.739	ng	98
57) 2,4-Dinitrotoluene	14.643	165	1200593	82.341	ng	98
58) Fluorene	15.319	166	3484502	76.824	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.901	232	970694	81.077	ng	100
60) Diethylphthalate	15.095	149	3799279	78.844	ng	98
61) 4-Chlorophenyl-phenylether	15.313	204	1725059	77.784	ng	97
62) 4-Nitroaniline	15.372	138	875375	89.199	ng	96
63) Azobenzene	15.613	77	3426736	77.538	ng	99
65) 4,6-Dinitro-2-methylph...	15.425	198	728771	86.490	ng	96
66) n-Nitrosodiphenylamine	15.543	169	3044661	76.603	ng	98
67) 4-Bromophenyl-phenylether	16.231	248	1157878	80.026	ng	99
68) Hexachlorobenzene	16.348	284	1395074	79.535	ng	99
69) Atrazine	16.525	200	1168753	81.114	ng	98
70) Pentachlorophenol	16.707	266	817045	90.003	ng	99
71) Phenanthrene	17.095	178	5340604	75.365	ng	99
72) Anthracene	17.189	178	5555906	77.415	ng	99
73) Carbazole	17.478	167	5166153	77.661	ng	99
74) Di-n-butylphthalate	18.066	149	6583572	79.879	ng	99
75) Fluoranthene	19.189	202	6351620	77.333	ng	100
77) Benzidine	19.389	184	2849177	68.307	ng	100
78) Pyrene	19.572	202	6497989	77.229	ng	99
80) Butylbenzylphthalate	20.525	149	3175233	82.388	ng	98
81) Benzo(a)anthracene	21.466	228	6696396	77.741	ng	99
82) 3,3'-Dichlorobenzidine	21.395	252	2698352	78.907	ng	99
83) Chrysene	21.536	228	6294743	77.125	ng	99
84) Bis(2-ethylhexyl)phtha...	21.407	149	4578538	82.917	ng	98
85) Di-n-octyl phthalate	22.636	149	7936082	81.539	ng	100
87) Indeno(1,2,3-cd)pyrene	28.424	276	9013339	78.751	ng	# 93
88) Benzo(b)fluoranthene	23.713	252	7112180	79.224	ng	99
89) Benzo(k)fluoranthene	23.783	252	7179421	78.564	ng	100
90) Benzo(a)pyrene	24.595	252	6983429	79.653	ng	99
91) Dibenzo(a,h)anthracene	28.501	278	7356566	78.965	ng	100
92) Benzo(g,h,i)perylene	29.554	276	7260298	78.545	ng	98

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060625\
 Data File : BP024867.D
 Acq On : 06 Jun 2025 15:18
 Operator : RC/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC080

Quant Time: Jun 06 16:05:27 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060625.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 06 15:53:04 2025
 Response via : Initial Calibration

