

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041425\
 Data File : BP024277.D
 Acq On : 14 Apr 2025 12:27
 Operator : RC/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC010

Quant Time: Apr 14 17:26:30 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP041425.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 14 17:03:26 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.734	152	237860	20.000	ng	0.00
21) Naphthalene-d8	10.510	136	1022895	20.000	ng	0.00
39) Acenaphthene-d10	14.363	164	658652	20.000	ng	0.00
64) Phenanthrene-d10	17.163	188	1310701	20.000	ng	0.00
76) Chrysene-d12	21.616	240	1408913	20.000	ng	0.00
86) Perylene-d12	24.980	264	1508235	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.352	112	272237	18.889	ng	0.00
7) Phenol-d6	6.910	99	369199	18.725	ng	0.00
23) Nitrobenzene-d5	8.881	82	345094	19.207	ng	0.00
42) 2,4,6-Tribromophenol	15.886	330	168790	18.715	ng	0.00
45) 2-Fluorobiphenyl	12.975	172	878366	20.247	ng	0.00
79) Terphenyl-d14	19.892	244	1405643	19.888	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.293	88	61433	9.984	ng	98
3) Pyridine	3.693	79	152243	9.506	ng	99
4) n-Nitrosodimethylamine	3.599	42	52702	9.868	ng	# 98
6) Aniline	7.069	93	176840	9.935	ng	99
8) 2-Chlorophenol	7.305	128	154194	9.573	ng	100
9) Benzaldehyde	6.887	77	111093m	10.879	ng	
10) Phenol	6.940	94	186113	9.401	ng	97
11) bis(2-Chloroethyl)ether	7.157	93	156684	9.794	ng	98
12) 1,3-Dichlorobenzene	7.622	146	176712	10.126	ng	97
13) 1,4-Dichlorobenzene	7.769	146	177067	10.006	ng	99
14) 1,2-Dichlorobenzene	8.081	146	171376	10.016	ng	99
15) Benzyl Alcohol	7.969	79	111176	8.797	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.252	45	177896	10.221	ng	98
17) 2-Methylphenol	8.175	107	119161	9.340	ng	98
18) Hexachloroethane	8.804	117	63912	10.007	ng	98
19) n-Nitroso-di-n-propyla...	8.528	70	119861	9.787	ng	100
20) 3+4-Methylphenols	8.499	107	162332	9.187	ng	99
22) Acetophenone	8.546	105	248823	9.787	ng	99
24) Nitrobenzene	8.922	77	170770	9.590	ng	98
25) Isophorone	9.446	82	307272	9.473	ng	99
26) 2-Nitrophenol	9.628	139	73320	8.575	ng	96
27) 2,4-Dimethylphenol	9.693	122	99186	9.152	ng	97
28) bis(2-Chloroethoxy)met...	9.922	93	217866	9.869	ng	99
29) 2,4-Dichlorophenol	10.163	162	136816	9.263	ng	99
30) 1,2,4-Trichlorobenzene	10.369	180	155965	9.760	ng	100
31) Naphthalene	10.557	128	538892	9.953	ng	99
32) Benzoic acid	9.781	122	94803m	8.095	ng	
33) 4-Chloroaniline	10.669	127	178451	9.433	ng	99
34) Hexachlorobutadiene	10.845	225	92662	9.885	ng	99
35) Caprolactam	11.440	113	47014	8.670	ng	98
36) 4-Chloro-3-methylphenol	11.798	107	158563	9.213	ng	99
37) 2-Methylnaphthalene	12.169	142	368122	9.810	ng	98
38) 1-Methylnaphthalene	12.387	142	365139	9.933	ng	100
40) 1,2,4,5-Tetrachloroben...	12.540	216	177188	9.789	ng	100
41) Hexachlorocyclopentadiene	12.516	237	60914	9.586	ng	96
43) 2,4,6-Trichlorophenol	12.781	196	110203	9.253	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041425\
 Data File : BP024277.D
 Acq On : 14 Apr 2025 12:27
 Operator : RC/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC010

Quant Time: Apr 14 17:26:30 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP041425.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 14 17:03:26 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.857	196	122104	9.251	ng	98
46) 1,1'-Biphenyl	13.187	154	490325	10.080	ng	100
47) 2-Chloronaphthalene	13.228	162	362419	9.985	ng	98
48) 2-Nitroaniline	13.434	65	88889	8.866	ng	95
49) Acenaphthylene	14.081	152	553215	9.711	ng	100
50) Dimethylphthalate	13.816	163	466314	9.704	ng	99
51) 2,6-Dinitrotoluene	13.934	165	95596	9.442	ng	96
52) Acenaphthene	14.428	154	357828	9.877	ng	99
53) 3-Nitroaniline	14.269	138	95237	9.067	ng	97
54) 2,4-Dinitrophenol	14.481	184	41132	7.588	ng	97
55) Dibenzofuran	14.769	168	594953	10.013	ng	99
56) 4-Nitrophenol	14.598	139	75995	8.416	ng	95
57) 2,4-Dinitrotoluene	14.734	165	121276	8.862	ng	# 95
58) Fluorene	15.428	166	459952	10.026	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.004	232	115645	9.470	ng	100
60) Diethylphthalate	15.198	149	488847	9.884	ng	100
61) 4-Chlorophenyl-phenyle...	15.428	204	222437	9.985	ng	97
62) 4-Nitroaniline	15.451	138	98638	8.857	ng	93
63) Azobenzene	15.722	77	455459	9.999	ng	96
65) 4,6-Dinitro-2-methylph...	15.516	198	62421	8.151	ng	96
66) n-Nitrosodiphenylamine	15.639	169	383558	9.759	ng	99
67) 4-Bromophenyl-phenylether	16.339	248	138724	9.703	ng	97
68) Hexachlorobenzene	16.457	284	161944	9.528	ng	98
69) Atrazine	16.628	200	109503	10.990	ng	98
70) Pentachlorophenol	16.810	266	101318	8.682	ng	96
71) Phenanthrene	17.210	178	711407	9.900	ng	99
72) Anthracene	17.304	178	666269	9.635	ng	100
73) Carbazole	17.586	167	654205	9.721	ng	99
74) Di-n-butylphthalate	18.180	149	806280	9.616	ng	99
75) Fluoranthene	19.292	202	835153	9.791	ng	98
77) Benzidine	19.498	184	99659	5.730	ng	99
78) Pyrene	19.669	202	870767	9.659	ng	99
80) Butylbenzylphthalate	20.627	149	343670	9.043	ng	99
81) Benzo(a)anthracene	21.598	228	854983	9.734	ng	98
82) 3,3'-Dichlorobenzidine	21.521	252	269772	8.857	ng	98
83) Chrysene	21.668	228	832263	9.887	ng	100
84) Bis(2-ethylhexyl)phtha...	21.539	149	517894	9.253	ng	99
85) Di-n-octyl phthalate	22.827	149	769654	8.584	ng	99
87) Indeno(1,2,3-cd)pyrene	28.792	276	983323	9.480	ng	99
88) Benzo(b)fluoranthene	23.915	252	885212	9.810	ng	99
89) Benzo(k)fluoranthene	23.986	252	852606	9.735	ng	99
90) Benzo(a)pyrene	24.815	252	732209	9.426	ng	99
91) Dibenzo(a,h)anthracene	28.862	278	821160	9.517	ng	99
92) Benzo(g,h,i)perylene	29.962	276	836993	9.530	ng	99

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

