

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090924\
 Data File : BP021863.D
 Acq On : 09 Sep 2024 12:59
 Operator : RC/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICCC040

Quant Time: Sep 10 00:02:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP090924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 09 23:51:05 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.758	152	161182	20.000	ng	0.00
21) Naphthalene-d8	10.534	136	568693	20.000	ng	0.00
39) Acenaphthene-d10	14.381	164	378217	20.000	ng	0.00
64) Phenanthrene-d10	17.169	188	847066	20.000	ng	0.00
76) Chrysene-d12	21.616	240	866364	20.000	ng	0.00
86) Perylene-d12	24.962	264	991783	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	1155844	97.661	ng	0.00
7) Phenol-d6	6.934	99	1645210	101.957	ng	0.00
23) Nitrobenzene-d5	8.905	82	1060582	98.319	ng	0.00
42) 2,4,6-Tribromophenol	15.886	330	384651	86.442	ng	0.00
45) 2-Fluorobiphenyl	12.993	172	1985006	85.533	ng	0.00
79) Terphenyl-d14	19.898	244	3559429	84.321	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.299	88	257070	42.145	ng	100
3) Pyridine	3.699	79	719793	45.348	ng	100
4) n-Nitrosodimethylamine	3.599	42	236150	42.292	ng	100
6) Aniline	7.093	93	747372	52.623	ng	100
8) 2-Chlorophenol	7.328	128	557368	50.369	ng	100
9) Benzaldehyde	6.905	77	440467	48.745	ng	100
10) Phenol	6.963	94	824797	50.977	ng	100
11) bis(2-Chloroethyl)ether	7.187	93	651966	51.715	ng	100
12) 1,3-Dichlorobenzene	7.652	146	511020	42.300	ng	100
13) 1,4-Dichlorobenzene	7.793	146	485377	40.145	ng	100
14) 1,2-Dichlorobenzene	8.110	146	468913	40.479	ng	100
15) Benzyl Alcohol	7.993	79	426908	40.373	ng	100
16) 2,2'-oxybis(1-Chloropr...	8.281	45	427640	38.996	ng	100
17) 2-Methylphenol	8.199	107	414298	40.976	ng	100
18) Hexachloroethane	8.834	117	196186	40.059	ng	100
19) n-Nitroso-di-n-propyla...	8.557	70	348318	39.598	ng	100
20) 3+4-Methylphenols	8.522	107	575937	41.237	ng	100
22) Acetophenone	8.569	105	802040	50.822	ng	100
24) Nitrobenzene	8.946	77	533635	48.783	ng	100
25) Isophorone	9.469	82	911139	43.066	ng	100
26) 2-Nitrophenol	9.652	139	186670	44.397	ng	100
27) 2,4-Dimethylphenol	9.716	122	260523	43.107	ng	100
28) bis(2-Chloroethoxy)met...	9.946	93	569943	41.826	ng	100
29) 2,4-Dichlorophenol	10.187	162	347881	44.184	ng	100
30) 1,2,4-Trichlorobenzene	10.399	180	382154	41.963	ng	100
31) Naphthalene	10.581	128	1220226	40.786	ng	100
32) Benzoic acid	9.840	122	292583	39.251	ng	100
33) 4-Chloroaniline	10.687	127	469354	42.484	ng	100
34) Hexachlorobutadiene	10.875	225	247873	41.874	ng	100
35) Caprolactam	11.475	113	144643	42.510	ng	100
36) 4-Chloro-3-methylphenol	11.816	107	504091	43.395	ng	100
37) 2-Methylnaphthalene	12.193	142	830778	42.597	ng	100
38) 1-Methylnaphthalene	12.416	142	823482	42.413	ng	100
40) 1,2,4,5-Tetrachloroben...	12.563	216	426563	42.950	ng	100
41) Hexachlorocyclopentadiene	12.551	237	148667	45.957	ng	100
43) 2,4,6-Trichlorophenol	12.804	196	289320	45.248	ng	100

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090924\
 Data File : BP021863.D
 Acq On : 09 Sep 2024 12:59
 Operator : RC/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICCC040

Quant Time: Sep 10 00:02:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP090924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 09 23:51:05 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.881	196	329147	45.184	ng	100
46) 1,1'-Biphenyl	13.210	154	1099309	43.091	ng	100
47) 2-Chloronaphthalene	13.251	162	855250	42.798	ng	100
48) 2-Nitroaniline	13.451	65	301603	44.865	ng	100
49) Acenaphthylene	14.098	152	1279210	42.142	ng	100
50) Dimethylphthalate	13.834	163	1125259	42.640	ng	100
51) 2,6-Dinitrotoluene	13.945	165	251510	43.181	ng	100
52) Acenaphthene	14.445	154	824822	41.898	ng	100
53) 3-Nitroaniline	14.281	138	262451	45.081	ng	100
54) 2,4-Dinitrophenol	14.492	184	127621	42.904	ng	100
55) Dibenzofuran	14.781	168	1299479	41.550	ng	100
56) 4-Nitrophenol	14.604	139	229869	44.735	ng	100
57) 2,4-Dinitrotoluene	14.745	165	361538	45.026	ng	100
58) Fluorene	15.439	166	1097688	42.543	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.016	232	284339	43.203	ng	100
60) Diethylphthalate	15.216	149	1204490	43.691	ng	100
61) 4-Chlorophenyl-phenyle...	15.434	204	528330	41.797	ng	100
62) 4-Nitroaniline	15.451	138	289970	45.190	ng	100
63) Azobenzene	15.734	77	1208342	41.172	ng	100
65) 4,6-Dinitro-2-methylph...	15.516	198	184117	44.324	ng	100
66) n-Nitrosodiphenylamine	15.651	169	933690	42.011	ng	100
67) 4-Bromophenyl-phenylether	16.345	248	329168	41.138	ng	100
68) Hexachlorobenzene	16.463	284	396811	41.193	ng	100
69) Atrazine	16.622	200	252667	36.047	ng	100
70) Pentachlorophenol	16.810	266	271806	44.012	ng	100
71) Phenanthrene	17.216	178	1731085	41.381	ng	100
72) Anthracene	17.310	178	1698559	42.310	ng	100
73) Carbazole	17.586	167	1737299	42.081	ng	100
74) Di-n-butylphthalate	18.175	149	2075958	44.077	ng	100
75) Fluoranthene	19.298	202	2198629	41.689	ng	100
77) Benzidine	19.498	184	759051	49.455	ng	100
78) Pyrene	19.675	202	2329123	41.662	ng	100
80) Butylbenzylphthalate	20.622	149	984449	43.526	ng	100
81) Benzo(a)anthracene	21.592	228	2278458	41.489	ng	100
82) 3,3'-Dichlorobenzidine	21.516	252	771484	43.286	ng	100
83) Chrysene	21.663	228	2163201	41.117	ng	100
84) Bis(2-ethylhexyl)phtha...	21.533	149	1452300	44.349	ng	100
85) Di-n-octyl phthalate	22.810	149	2417575	45.314	ng	100
87) Indeno(1,2,3-cd)pyrene	28.745	276	2590148m	41.398	ng	
88) Benzo(b)fluoranthene	23.898	252	2380527	41.409	ng	100
89) Benzo(k)fluoranthene	23.963	252	2289625	41.750	ng	100
90) Benzo(a)pyrene	24.786	252	2073382	41.965	ng	100
91) Dibenzo(a,h)anthracene	28.821	278	2114709	41.188	ng	100
92) Benzo(g,h,i)perylene	29.927	276	2245482	41.721	ng	100

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

