

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011521\
 Data File : BP004580.D
 Acq On : 15 Jan 2021 17:06
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Quant Time: Jan 15 18:04:45 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP011221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 12 15:19:32 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	301481	20.00	ng	-0.01
21) Naphthalene-d8	10.599	136	1064332	20.00	ng	#-0.01
39) Acenaphthene-d10	14.457	164	621301	20.00	ng	-0.01
64) Phenanthrene-d10	17.216	188	1259521	20.00	ng	-0.01
76) Chrysene-d12	21.310	240	1187985	20.00	ng	-0.01
86) Perylene-d12	23.663	264	1250269	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	1355159	78.35	ng	0.00
7) Phenol-d6	6.981	99	1741204	73.50	ng	0.00
23) Nitrobenzene-d5	8.963	82	1420591	75.46	ng	-0.01
42) 2,4,6-Tribromophenol	15.957	330	718611	66.52	ng	0.00
45) 2-Fluorobiphenyl	13.075	172	3699247	77.40	ng	0.00
79) Terphenyl-d14	19.781	244	4876758	75.00	ng	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.323	88	293597	41.99	ng	93
3) Pyridine	3.711	79	750980	40.08	ng	95
4) n-Nitrosodimethylamine	3.628	42	234335	39.56	ng	92
6) Aniline	7.134	93	969653	35.98	ng	96
8) 2-Chlorophenol	7.375	128	731333	37.23	ng	95
9) Benzaldehyde	6.946	77	412713	40.77	ng	92
10) Phenol	7.011	94	837777	36.88	ng	99
11) bis(2-Chloroethyl)ether	7.234	93	674646	37.11	ng	96
12) 1,3-Dichlorobenzene	7.693	146	901391	37.11	ng	98
13) 1,4-Dichlorobenzene	7.840	146	910288	37.05	ng	99
14) 1,2-Dichlorobenzene	8.158	146	862210	36.75	ng	100
15) Benzyl Alcohol	8.046	79	535700	34.90	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.334	45	703237	36.48	ng	94
17) 2-Methylphenol	8.252	107	575439	35.90	ng	99
18) Hexachloroethane	8.881	117	309942	36.50	ng	95
19) n-Nitroso-di-n-propyla...	8.611	70	440004	33.11	ng	94
20) 3+4-Methylphenols	8.581	107	775200	35.45	ng	99
22) Acetophenone	8.628	105	992259	37.77	ng	# 97
24) Nitrobenzene	9.011	77	698483	37.92	ng	92
25) Isophorone	9.534	82	1220039	35.94	ng	96
26) 2-Nitrophenol	9.716	139	397415	39.14	ng	94
27) 2,4-Dimethylphenol	9.781	122	533300	37.92	ng	98
28) bis(2-Chloroethoxy)met...	10.010	93	857348	36.80	ng	99
29) 2,4-Dichlorophenol	10.258	162	699317	37.44	ng	99
30) 1,2,4-Trichlorobenzene	10.469	180	837295	37.27	ng	98
31) Naphthalene	10.652	128	2190038	37.46	ng	100
32) Benzoic acid	9.952	122	406723	37.31	ng	91
33) 4-Chloroaniline	10.763	127	905311	35.86	ng	98
34) Hexachlorobutadiene	10.934	225	547716	36.28	ng	100
35) Caprolactam	11.546	113	194391	32.60	ng	# 75
36) 4-Chloro-3-methylphenol	11.904	107	622792	35.02	ng	99
37) 2-Methylnaphthalene	12.269	142	1537716	35.78	ng	100
38) 1-Methylnaphthalene	12.487	142	1455967	35.69	ng	98
40) 1,2,4,5-Tetrachloroben...	12.640	216	881370	39.49	ng	99
41) Hexachlorocyclopentadiene	12.616	237	326376	38.09	ng	99
43) 2,4,6-Trichlorophenol	12.887	196	566555	39.00	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.963	196	582534	38.59	ng	99
46) 1,1'-Biphenyl	13.287	154	1930952	38.88	ng	99
47) 2-Chloronaphthalene	13.328	162	1604160	39.02	ng	98
48) 2-Nitroaniline	13.534	65	362229	38.11	ng	86
49) Acenaphthylene	14.181	152	2327067	37.89	ng	99
50) Dimethylphthalate	13.916	163	1861155	36.15	ng	99
51) 2,6-Dinitrotoluene	14.040	165	408858	37.25	ng	99
52) Acenaphthene	14.522	154	1409596	37.65	ng	99
53) 3-Nitroaniline	14.369	138	434352	36.36	ng	94
54) 2,4-Dinitrophenol	14.593	184	188388	36.23	ng	94
55) Dibenzofuran	14.857	168	2252852	37.16	ng	97
56) 4-Nitrophenol	14.698	139	296303	35.73	ng	92
57) 2,4-Dinitrotoluene	14.834	165	547294	36.47	ng	93
58) Fluorene	15.510	166	1774468	36.66	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.093	232	509044	35.93	ng	# 91
60) Diethylphthalate	15.281	149	1764317	35.04	ng	98
61) 4-Chlorophenyl-phenylether	15.504	204	986316	36.06	ng	92
62) 4-Nitroaniline	15.540	138	451042	35.75	ng	97
63) Azobenzene	15.798	77	1386095	36.36	ng	96
65) 4,6-Dinitro-2-methylph...	15.610	198	280084	38.18	ng	96
66) n-Nitrosodiphenylamine	15.722	169	1521710	38.97	ng	100
67) 4-Bromophenyl-phenylether	16.404	248	639791	38.01	ng	94
68) Hexachlorobenzene	16.528	284	742275	36.91	ng	96
69) Atrazine	16.681	200	482092	35.10	ng	98
70) Pentachlorophenol	16.875	266	359330	39.82	ng	99
71) Phenanthrene	17.263	178	2728186	37.78	ng	100
72) Anthracene	17.351	178	2652699	37.91	ng	100
73) Carbazole	17.628	167	2412178	36.99	ng	99
74) Di-n-butylphthalate	18.175	149	2849606	36.15	ng	99
75) Fluoranthene	19.234	202	3103179	35.32	ng	100
77) Benzidine	19.416	184	866145	27.87	ng	100
78) Pyrene	19.586	202	3157787	39.07	ng	100
80) Butylbenzylphthalate	20.457	149	1241740	37.98	ng	95
81) Benzo(a)anthracene	21.298	228	3029290	37.49	ng	100
82) 3,3'-Dichlorobenzidine	21.228	252	1074528	36.89	ng	99
83) Chrysene	21.351	228	2892006	37.60	ng	99
84) Bis(2-ethylhexyl)phtha...	21.216	149	1779128	38.11	ng	98
85) Di-n-octyl phthalate	22.122	149	2979765	37.59	ng	# 97
87) Indeno(1,2,3-cd)pyrene	26.086	276	3829154	39.69	ng	98
88) Benzo(b)fluoranthene	22.951	252	3216702	39.75	ng	99
89) Benzo(k)fluoranthene	22.998	252	2968351	38.46	ng	99
90) Benzo(a)pyrene	23.557	252	2954767	39.14	ng	99
91) Dibenzo(a,h)anthracene	26.092	278	3144940	39.50	ng	98
92) Benzo(g,h,i)perylene	26.827	276	3091741	39.43	ng	97

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

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