

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091322\
 Data File : BP011685.D
 Acq On : 13 Sep 2022 14:21
 Operator : CG/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 09/14/2022
 Supervised By : Jagrut Upadhyay 09/14/2022

Quant Time: Sep 13 15:11:08 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 13 14:11:58 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	409630	20.000	ng	0.00
21) Naphthalene-d8	10.769	136	1538962	20.000	ng	# 0.00
39) Acenaphthene-d10	14.592	164	888074	20.000	ng	0.00
64) Phenanthrene-d10	17.351	188	1764124	20.000	ng	# 0.00
76) Chrysene-d12	21.433	240	1642517	20.000	ng	# 0.00
86) Perylene-d12	23.892	264	1621677	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.511	112	3594775	155.206	ng	0.00
7) Phenol-d6	7.116	99	4681147	152.743	ng	0.00
23) Nitrobenzene-d5	9.122	82	4196846	167.397	ng	0.00
42) 2,4,6-Tribromophenol	16.086	330	1303322	129.892	ng	0.00
45) 2-Fluorobiphenyl	13.222	172	8672484	139.647	ng	0.00
79) Terphenyl-d14	19.904	244	11237087	143.292	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.393	88	761363	83.526	ng	# 79
3) Pyridine	3.799	79	2276150m	91.702	ng	
4) n-Nitrosodimethylamine	3.705	42	800408	77.214	ng	# 36
6) Aniline	7.281	93	2887892	82.024	ng	89
8) 2-Chlorophenol	7.516	128	2067003	78.049	ng	90
9) Benzaldehyde	7.093	77	1313772m	73.025	ng	
10) Phenol	7.140	94	2620949	81.780	ng	81
11) bis(2-Chloroethyl)ether	7.375	93	1969573	78.193	ng	91
12) 1,3-Dichlorobenzene	7.840	146	2223578	73.397	ng	# 93
13) 1,4-Dichlorobenzene	7.993	146	2270659	73.521	ng	97
14) 1,2-Dichlorobenzene	8.310	146	2144207	73.093	ng	97
15) Benzyl Alcohol	8.199	79	1684379	82.176	ng	# 88
16) 2,2'-oxybis(1-Chloropr...	8.487	45	2393136	73.713	ng	90
17) 2-Methylphenol	8.399	107	1705738	80.521	ng	# 91
18) Hexachloroethane	9.040	117	788120	75.789	ng	96
19) n-Nitroso-di-n-propyla...	8.775	70	1442231	80.809	ng	# 78
20) 3+4-Methylphenols	8.728	107	2354038	81.994	ng	92
22) Acetophenone	8.781	105	2895373	77.835	ng	# 85
24) Nitrobenzene	9.163	77	2158916	83.389	ng	# 87
25) Isophorone	9.699	82	3756093	83.098	ng	# 94
26) 2-Nitrophenol	9.875	139	1028163	105.906	ng	# 79
27) 2,4-Dimethylphenol	9.934	122	1613136	82.307	ng	89
28) bis(2-Chloroethoxy)met...	10.175	93	2296919	81.235	ng	98
29) 2,4-Dichlorophenol	10.410	162	1817207	79.513	ng	97
30) 1,2,4-Trichlorobenzene	10.628	180	1902532	70.943	ng	98
31) Naphthalene	10.816	128	6312475	76.703	ng	98
32) Benzoic acid	10.099	122	1376784	119.934	ng	# 87
33) 4-Chloroaniline	10.928	127	2645257	81.477	ng	# 91
34) Hexachlorobutadiene	11.104	225	1062214	66.570	ng	98
35) Caprolactam	11.740	113	609234	92.518	ng	# 64
36) 4-Chloro-3-methylphenol	12.045	107	1931238	85.759	ng	91
37) 2-Methylnaphthalene	12.422	142	4265388	76.131	ng	97
38) 1-Methylnaphthalene	12.640	142	4148397	76.016	ng	99
40) 1,2,4,5-Tetrachloroben...	12.787	216	1913364	68.597	ng	98
41) Hexachlorocyclopentadiene	12.769	237	989628	75.050	ng	99
43) 2,4,6-Trichlorophenol	13.028	196	1387104	83.099	ng	97

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091322\
 Data File : BP011685.D
 Acq On : 13 Sep 2022 14:21
 Operator : CG/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 09/14/2022
 Supervised By : Jagrut Upadhyay 09/14/2022

Quant Time: Sep 13 15:11:08 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 13 14:11:58 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.098	196	1535612	80.589	ng	97
46) 1,1'-Biphenyl	13.434	154	5068375	74.301	ng	100
47) 2-Chloronaphthalene	13.475	162	4039990	76.192	ng	99
48) 2-Nitroaniline	13.675	65	1244994	102.257	ng	# 77
49) Acenaphthylene	14.316	152	6434110	79.089	ng	99
50) Dimethylphthalate	14.057	163	4968195	76.842	ng	100
51) 2,6-Dinitrotoluene	14.169	165	1100557	88.069	ng	# 83
52) Acenaphthene	14.663	154	4219321m	80.111	ng	
53) 3-Nitroaniline	14.504	138	1313147	94.744	ng	86
54) 2,4-Dinitrophenol	14.704	184	580257	139.177	ng	# 81
55) Dibenzofuran	14.992	168	5979586	75.353	ng	98
56) 4-Nitrophenol	14.798	139	1092259	97.971	ng	# 71
57) 2,4-Dinitrotoluene	14.957	165	1488992	94.879	ng	# 81
58) Fluorene	15.645	166	4801025	73.885	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.216	232	1273787	79.208	ng	97
60) Diethylphthalate	15.422	149	4979483	79.205	ng	99
61) 4-Chlorophenyl-phenylether	15.639	204	2251032	70.262	ng	95
62) 4-Nitroaniline	15.669	138	1369797	97.337	ng	# 77
63) Azobenzene	15.933	77	4712339	84.308	ng	87
65) 4,6-Dinitro-2-methylph...	15.722	198	799086	117.841	ng	# 84
66) n-Nitrosodiphenylamine	15.851	169	4194061	78.158	ng	97
67) 4-Bromophenyl-phenylether	16.533	248	1368225	71.362	ng	98
68) Hexachlorobenzene	16.651	284	1492886	65.984	ng	97
69) Atrazine	16.810	200	1368444	85.786	ng	96
70) Pentachlorophenol	16.992	266	1009160	82.049	ng	99
71) Phenanthrene	17.392	178	7564562	76.847	ng	97
72) Anthracene	17.486	178	7370088	79.034	ng	97
73) Carbazole	17.751	167	7071804	81.967	ng	98
74) Di-n-butylphthalate	18.304	149	8285784	85.076	ng	98
75) Fluoranthene	19.351	202	8489991	78.237	ng	93
77) Benzidine	19.527	184	2617264	83.365	ng	98
78) Pyrene	19.704	202	8644990	79.184	ng	96
80) Butylbenzylphthalate	20.574	149	3891709	101.911	ng	93
81) Benzo(a)anthracene	21.416	228	8395510	77.865	ng	94
82) 3,3'-Dichlorobenzidine	21.345	252	2752500	78.257	ng	99
83) Chrysene	21.474	228	7941404	77.841	ng	96
84) Bis(2-ethylhexyl)phtha...	21.333	149	5340644	90.820	ng	98
85) Di-n-octyl phthalate	22.286	149	9215549	93.201	ng	# 90
87) Indeno(1,2,3-cd)pyrene	26.492	276	9888644	80.137	ng	# 92
88) Benzo(b)fluoranthene	23.145	252	8274315	81.331	ng	# 97
89) Benzo(k)fluoranthene	23.198	252	8261370	78.715	ng	# 96
90) Benzo(a)pyrene	23.786	252	7017761	83.484	ng	# 95
91) Dibenzo(a,h)anthracene	26.515	278	8092439	78.175	ng	# 95
92) Benzo(g,h,i)perylene	27.286	276	8061719	81.006	ng	# 92

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

