

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071221\
 Data File : BP006183.D
 Acq On : 12 Jul 2021 14:18
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 7/13/2021 1:42:09 PM

Quant Time: Jul 12 15:00:10 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 12 14:59:51 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	210589	20.000	ng	0.00
21) Naphthalene-d8	10.663	136	799010	20.000	ng	0.00
39) Acenaphthene-d10	14.498	164	423968	20.000	ng	0.00
64) Phenanthrene-d10	17.251	188	756519	20.000	ng	# 0.00
76) Chrysene-d12	21.327	240	748641	20.000	ng	# 0.00
86) Perylene-d12	23.715	264	903775	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	972411	79.452	ng	0.00
7) Phenol-d6	7.045	99	1265244	78.155	ng	0.00
23) Nitrobenzene-d5	9.034	82	1011258	79.944	ng	0.00
42) 2,4,6-Tribromophenol	15.998	330	432630	85.710	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	2486158	86.690	ng	0.00
79) Terphenyl-d14	19.804	244	3122412	82.232	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.304	88	182229	37.948	ng	93
3) Pyridine	3.728	79	490506	40.200	ng	90
4) n-Nitrosodimethylamine	3.634	42	179257	37.351	ng	# 83
6) Aniline	7.192	93	657418	38.489	ng	98
8) 2-Chlorophenol	7.434	128	567785	39.611	ng	96
9) Benzaldehyde	7.004	77	331851	37.559	ng	92
10) Phenol	7.075	94	611523	39.331	ng	91
11) bis(2-Chloroethyl)ether	7.292	93	460608	39.205	ng	99
12) 1,3-Dichlorobenzene	7.745	146	612383	39.871	ng	# 95
13) 1,4-Dichlorobenzene	7.898	146	622920	39.712	ng	98
14) 1,2-Dichlorobenzene	8.210	146	591845	39.633	ng	98
15) Benzyl Alcohol	8.116	79	378059m	38.628	ng	
16) 2,2'-oxybis(1-Chloropr...	8.398	45	515832	36.968	ng	90
17) 2-Methylphenol	8.322	107	417515	39.474	ng	94
18) Hexachloroethane	8.934	117	206984	39.637	ng	96
19) n-Nitroso-di-n-propyla...	8.675	70	336603	38.139	ng	97
20) 3+4-Methylphenols	8.657	107	569242	39.836	ng	93
22) Acetophenone	8.692	105	728465	40.056	ng	98
24) Nitrobenzene	9.075	77	510080	39.871	ng	89
25) Isophorone	9.604	82	940210	39.154	ng	95
26) 2-Nitrophenol	9.786	139	304762	42.852	ng	# 84
27) 2,4-Dimethylphenol	9.851	122	431327	41.085	ng	95
28) bis(2-Chloroethoxy)met...	10.081	93	530511	39.865	ng	100
29) 2,4-Dichlorophenol	10.334	162	489487	42.434	ng	100
30) 1,2,4-Trichlorobenzene	10.528	180	510402	41.559	ng	98
31) Naphthalene	10.716	128	1657237	39.787	ng	99
32) Benzoic acid	10.063	122	320155	41.284	ng	89
33) 4-Chloroaniline	10.839	127	657510	40.893	ng	97
34) Hexachlorobutadiene	10.992	225	288887	42.721	ng	97
35) Caprolactam	11.657	113	146884	40.394	ng	95
36) 4-Chloro-3-methylphenol	11.980	107	494525	39.701	ng	96
37) 2-Methylnaphthalene	12.328	142	1148351	39.736	ng	96
38) 1-Methylnaphthalene	12.545	142	1077998	39.414	ng	# 99
40) 1,2,4,5-Tetrachloroben...	12.698	216	475605	45.223	ng	98
41) Hexachlorocyclopentadiene	12.669	237	60773	38.189	ng	97
43) 2,4,6-Trichlorophenol	12.945	196	347783	45.709	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.033	196	372728	45.659	ng	# 97
46) 1,1'-Biphenyl	13.333	154	1370066	43.044	ng	99
47) 2-Chloronaphthalene	13.380	162	1060083	43.293	ng	99
48) 2-Nitroaniline	13.598	65	259809	42.698	ng	# 82
49) Acenaphthylene	14.222	152	1625734	41.232	ng	100
50) Dimethylphthalate	13.963	163	1270220	41.664	ng	99
51) 2,6-Dinitrotoluene	14.092	165	292812	42.601	ng	89
52) Acenaphthene	14.563	154	945355	39.415	ng	98
53) 3-Nitroaniline	14.422	138	295724	42.467	ng	94
54) 2,4-Dinitrophenol	14.657	184	127811	40.348	ng	97
55) Dibenzofuran	14.898	168	1462229	39.163	ng	98
56) 4-Nitrophenol	14.769	139	183711	38.310	ng	# 82
57) 2,4-Dinitrotoluene	14.886	165	380078	41.058	ng	95
58) Fluorene	15.551	166	1156347	38.550	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.139	232	267585	41.892	ng	# 80
60) Diethylphthalate	15.321	149	1209226	38.676	ng	99
61) 4-Chlorophenyl-phenyle...	15.539	204	515719	39.219	ng	93
62) 4-Nitroaniline	15.592	138	267649	38.263	ng	# 81
63) Azobenzene	15.833	77	981588	37.606	ng	96
65) 4,6-Dinitro-2-methylph...	15.657	198	199432	41.078	ng	89
66) n-Nitrosodiphenylamine	15.757	169	991262	40.241	ng	99
67) 4-Bromophenyl-phenylether	16.433	248	316217	42.401	ng	91
68) Hexachlorobenzene	16.557	284	377211	42.611	ng	95
69) Atrazine	16.716	200	321913	41.626	ng	95
70) Pentachlorophenol	16.916	266	187681	45.937	ng	96
71) Phenanthrene	17.292	178	1722360	40.313	ng	99
72) Anthracene	17.386	178	1698858	40.588	ng	100
73) Carbazole	17.663	167	1708247	41.186	ng	99
74) Di-n-butylphthalate	18.198	149	2092500	41.913	ng	99
75) Fluoranthene	19.257	202	1926738	41.146	ng	96
77) Benzidine	19.445	184	882929	39.773	ng	99
78) Pyrene	19.609	202	1982850	39.186	ng	99
80) Butylbenzylphthalate	20.474	149	1003091	39.965	ng	90
81) Benzo(a)anthracene	21.315	228	1916175	40.469	ng	99
82) 3,3'-Dichlorobenzidine	21.245	252	591480	42.583	ng	# 98
83) Chrysene	21.368	228	1869818	39.959	ng	99
84) Bis(2-ethylhexyl)phtha...	21.227	149	1499784	40.484	ng	99
85) Di-n-octyl phthalate	22.133	149	2706284	41.476	ng	100
87) Indeno(1,2,3-cd)pyrene	26.233	276	2743088	41.168	ng	# 91
88) Benzo(b)fluoranthene	22.986	252	2227339	40.628	ng	# 97
89) Benzo(k)fluoranthene	23.033	252	2116305	39.494	ng	# 96
90) Benzo(a)pyrene	23.609	252	2069440	40.226	ng	# 96
91) Dibenzo(a,h)anthracene	26.239	278	2382708	41.529	ng	# 94
92) Benzo(g,h,i)perylene	27.003	276	2308131	40.877	ng	# 91

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

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