

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM021423\
 Data File : BM038724.D
 Acq On : 14 Feb 2023 11:00
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Christian
 Giraldo

Quant Time: Feb 14 15:50:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM020123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 14 15:49:49 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.898	152	51792	20.000	ng	0.00
21) Naphthalene-d8	10.710	136	174812	20.000	ng	0.00
39) Acenaphthene-d10	14.545	164	112108	20.000	ng	0.00
64) Phenanthrene-d10	17.286	188	225672	20.000	ng	# 0.00
76) Chrysene-d12	21.462	240	194643	20.000	ng	0.00
86) Perylene-d12	23.838	264	244681	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.457	112	211156	79.865	ng	0.00
7) Phenol-d6	7.081	99	249223	82.648	ng	0.00
23) Nitrobenzene-d5	9.081	82	244396	79.410	ng	0.00
42) 2,4,6-Tribromophenol	16.033	330	125661	82.466	ng	0.00
45) 2-Fluorobiphenyl	13.168	172	732302	83.160	ng	0.00
79) Terphenyl-d14	19.903	244	948226	85.968	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.310	88	42840	36.754	ng	94
3) Pyridine	3.728	79	109888	39.647	ng	97
4) n-Nitrosodimethylamine	3.640	42	47545	37.998	ng	# 94
6) Aniline	7.234	93	154217	40.999	ng	95
8) 2-Chlorophenol	7.463	128	117274	40.650	ng	97
9) Benzaldehyde	7.045	77	49921	37.561	ng	89
10) Phenol	7.104	94	123138	40.421	ng	91
11) bis(2-Chloroethyl)ether	7.334	93	103346	41.890	ng	97
12) 1,3-Dichlorobenzene	7.786	146	149264	40.494	ng	98
13) 1,4-Dichlorobenzene	7.939	146	147857	39.746	ng	100
14) 1,2-Dichlorobenzene	8.251	146	142173	39.893	ng	98
15) Benzyl Alcohol	8.157	79	84419	38.036	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.434	45	89700	38.385	ng	98
17) 2-Methylphenol	8.357	107	96010	41.088	ng	98
18) Hexachloroethane	8.975	117	48505	40.374	ng	97
19) n-Nitroso-di-n-propyla...	8.722	70	73650	40.520	ng	96
20) 3+4-Methylphenols	8.686	107	129002	41.407	ng	96
22) Acetophenone	8.739	105	162562	38.994	ng	# 98
24) Nitrobenzene	9.122	77	120855	38.688	ng	94
25) Isophorone	9.645	82	206192	38.244	ng	98
26) 2-Nitrophenol	9.828	139	72804	41.883	ng	96
27) 2,4-Dimethylphenol	9.892	122	100999	40.271	ng	97
28) bis(2-Chloroethoxy)met...	10.128	93	132641	40.257	ng	97
29) 2,4-Dichlorophenol	10.363	162	136756	38.381	ng	98
30) 1,2,4-Trichlorobenzene	10.569	180	138533	36.551	ng	98
31) Naphthalene	10.763	128	367421	40.958	ng	99
32) Benzoic acid	10.051	122	76395	37.630	ng	97
33) 4-Chloroaniline	10.880	127	160014	40.701	ng	94
34) Hexachlorobutadiene	11.033	225	68774	36.223	ng	100
35) Caprolactam	11.686	113	30055	39.166	ng	99
36) 4-Chloro-3-methylphenol	12.010	107	101670	38.589	ng	95
37) 2-Methylnaphthalene	12.369	142	290891	40.127	ng	98
38) 1-Methylnaphthalene	12.586	142	272773	39.933	ng	98
40) 1,2,4,5-Tetrachloroben...	12.733	216	117631	40.317	ng	99
41) Hexachlorocyclopentadiene	12.704	237	65112	35.964	ng	97
43) 2,4,6-Trichlorophenol	12.980	196	94429	41.197	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	13.057	196	107998	40.039	ng	95	02/15/2023
46) 1,1'-Biphenyl	13.380	154	389290	43.836	ng	100	Supervised By :Jagrut Upadhyay
47) 2-Chloronaphthalene	13.421	162	296589	42.443	ng	98	
48) 2-Nitroaniline	13.639	65	62178	39.814	ng	94	
49) Acenaphthylene	14.268	152	485721	44.130	ng	99	
50) Dimethylphthalate	14.010	163	375691	41.065	ng	100	
51) 2,6-Dinitrotoluene	14.133	165	79898	42.969	ng	95	02/17/2023
52) Acenaphthene	14.610	154	283236	43.097	ng	99	
53) 3-Nitroaniline	14.468	138	82626	42.645	ng	97	
54) 2,4-Dinitrophenol	14.674	184	48465	40.044	ng	99	
55) Dibenzofuran	14.939	168	472924	41.788	ng	99	
56) 4-Nitrophenol	14.780	139	64370	43.164	ng	93	
57) 2,4-Dinitrotoluene	14.921	165	108770	39.927	ng	# 98	
58) Fluorene	15.592	166	377360	42.141	ng	98	
59) 2,3,4,6-Tetrachlorophenol	15.168	232	76982	41.006	ng	97	
60) Diethylphthalate	15.362	149	375896	43.371	ng	97	
61) 4-Chlorophenyl-phenyle...	15.580	204	155999	41.064	ng	99	
62) 4-Nitroaniline	15.627	138	83023m	41.807	ng		
63) Azobenzene	15.874	77	296872	44.521	ng	97	
65) 4,6-Dinitro-2-methylph...	15.680	198	59989	42.226	ng	96	
66) n-Nitrosodiphenylamine	15.798	169	315499	42.817	ng	98	
67) 4-Bromophenyl-phenylether	16.474	248	91835	42.145	ng	95	
68) Hexachlorobenzene	16.586	284	111987	42.052	ng	98	
69) Atrazine	16.751	200	89185	41.760	ng	100	
70) Pentachlorophenol	16.939	266	80436	41.434	ng	98	
71) Phenanthrene	17.327	178	561356	41.893	ng	99	
72) Anthracene	17.421	178	575307	41.968	ng	99	
73) Carbazole	17.698	167	544802	42.614	ng	100	
74) Di-n-butylphthalate	18.245	149	675613	46.775	ng	99	
75) Fluoranthene	19.339	202	592957	42.146	ng	98	
77) Benzidine	19.533	184	190527	33.581	ng	98	
78) Pyrene	19.703	202	626424	41.802	ng	99	
80) Butylbenzylphthalate	20.592	149	312109	42.085	ng	94	
81) Benzo(a)anthracene	21.444	228	576368	41.862	ng	99	
82) 3,3'-Dichlorobenzidine	21.386	252	202034	42.652	ng	98	
83) Chrysene	21.503	228	546967	41.231	ng	98	
84) Bis(2-ethylhexyl)phtha...	21.362	149	460725	42.614	ng	98	
85) Di-n-octyl phthalate	22.274	149	798154	45.248	ng	100	
87) Indeno(1,2,3-cd)pyrene	26.315	276	809580	42.007	ng	97	
88) Benzo(b)fluoranthene	23.115	252	610495	41.657	ng	98	
89) Benzo(k)fluoranthene	23.162	252	626721	42.112	ng	# 98	
90) Benzo(a)pyrene	23.738	252	606469	41.846	ng	98	
91) Dibenzo(a,h)anthracene	26.326	278	678080	41.819	ng	98	
92) Benzo(g,h,i)perylene	27.079	276	666964	40.692	ng	97	

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

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