

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102622\
 Data File : BF130781.D
 Acq On : 26 Oct 2022 11:53
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 10/27/2022
 Supervised By : Jagrut Upadhyay 11/01/2022

Quant Time: Oct 27 09:01:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 27 08:55:38 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.563	152	59454	20.000	ng	0.00
21) Naphthalene-d8	7.845	136	225483	20.000	ng	0.00
39) Acenaphthene-d10	9.592	164	123305	20.000	ng	0.00
64) Phenanthrene-d10	11.075	188	216018	20.000	ng	0.00
76) Chrysene-d12	13.716	240	139242	20.000	ng	# 0.00
86) Perylene-d12	15.098	264	112393	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.175	112	266997	77.440	ng	0.00
7) Phenol-d6	6.228	99	329623	79.017	ng	0.00
23) Nitrobenzene-d5	7.140	82	318446	74.918	ng	0.00
42) 2,4,6-Tribromophenol	10.386	330	95423	81.376	ng	0.00
45) 2-Fluorobiphenyl	8.922	172	650269	80.367	ng	0.00
79) Terphenyl-d14	12.657	244	695101	84.125	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.122	88	54524	26.580	ng	99
3) Pyridine	2.787	79	136282	34.461	ng	99
4) n-Nitrosodimethylamine	2.757	42	62747	35.713	ng	94
6) Aniline	6.234	93	198935	41.392	ng	# 82
8) 2-Chlorophenol	6.357	128	158054	39.618	ng	96
9) Benzaldehyde	6.122	77	89895	36.909	ng	96
10) Phenol	6.240	94	186734	40.603	ng	87
11) bis(2-Chloroethyl)ether	6.310	93	134040	39.240	ng	99
12) 1,3-Dichlorobenzene	6.504	146	169795	39.179	ng	100
13) 1,4-Dichlorobenzene	6.581	146	177492	39.818	ng	99
14) 1,2-Dichlorobenzene	6.734	146	163995	38.940	ng	99
15) Benzyl Alcohol	6.734	79	119442	39.034	ng	98
16) 2,2'-oxybis(1-Chloropr...	6.845	45	164321	37.502	ng	# 64
17) 2-Methylphenol	6.845	107	121400	36.990	ng	96
18) Hexachloroethane	7.063	117	63047	37.870	ng	93
19) n-Nitroso-di-n-propyla...	6.992	70	102198	36.256	ng	97
20) 3+4-Methylphenols	7.004	107	163818	38.083	ng	93
22) Acetophenone	6.992	105	211492	38.450	ng	# 96
24) Nitrobenzene	7.163	77	158900	36.749	ng	94
25) Isophorone	7.398	82	260309	37.170	ng	99
26) 2-Nitrophenol	7.475	139	81336	37.518	ng	96
27) 2,4-Dimethylphenol	7.522	122	113966	38.854	ng	99
28) bis(2-Chloroethoxy)met...	7.610	93	150286	37.046	ng	99
29) 2,4-Dichlorophenol	7.728	162	129117	39.921	ng	98
30) 1,2,4-Trichlorobenzene	7.792	180	140603	41.396	ng	97
31) Naphthalene	7.869	128	464437	39.948	ng	100
32) Benzoic acid	7.704	122	50923m	32.246	ng	
33) 4-Chloroaniline	7.934	127	185270	41.421	ng	100
34) Hexachlorobutadiene	7.981	225	86263	39.947	ng	98
35) Caprolactam	8.322	113	40267	39.835	ng	92
36) 4-Chloro-3-methylphenol	8.434	107	139195	39.820	ng	96
37) 2-Methylnaphthalene	8.557	142	304127	39.925	ng	99
38) 1-Methylnaphthalene	8.657	142	293321	38.838	ng	100
40) 1,2,4,5-Tetrachloroben...	8.728	216	133239	40.722	ng	98
41) Hexachlorocyclopentadiene	8.704	237	2203	20.129	ng	98
43) 2,4,6-Trichlorophenol	8.851	196	81933	39.787	ng	94

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44) 2,4,5-Trichlorophenol	8.904	196	83998	37.507	ng	97
46) 1,1'-Biphenyl	9.022	154	354143	41.177	ng	99
47) 2-Chloronaphthalene	9.045	162	291815	40.801	ng	99
48) 2-Nitroaniline	9.163	65	88263	40.206	ng	98
49) Acenaphthylene	9.457	152	439586	41.713	ng	99
50) Dimethylphthalate	9.333	163	342528	41.299	ng	100
51) 2,6-Dinitrotoluene	9.404	165	76239	41.048	ng	97
52) Acenaphthene	9.628	154	267164	40.787	ng	99
53) 3-Nitroaniline	9.575	138	84110	39.743	ng	98
54) 2,4-Dinitrophenol	9.722	184	15878m	27.384	ng	
55) Dibenzofuran	9.804	168	413838	41.046	ng	96
56) 4-Nitrophenol	9.786	139	2168m	10.303	ng	
57) 2,4-Dinitrotoluene	9.810	165	102841	41.105	ng	# 87
58) Fluorene	10.145	166	343081	40.915	ng	99
59) 2,3,4,6-Tetrachlorophenol	9.933	232	67920	38.147	ng	96
60) Diethylphthalate	10.028	149	332263	38.828	ng	99
61) 4-Chlorophenyl-phenyle...	10.133	204	155307	41.324	ng	92
62) 4-Nitroaniline	10.192	138	77530	38.620	ng	99
63) Azobenzene	10.298	77	302348	40.241	ng	99
65) 4,6-Dinitro-2-methylph...	10.228	198	52602	40.743	ng	99
66) n-Nitrosodiphenylamine	10.263	169	283885	38.122	ng	98
67) 4-Bromophenyl-phenylether	10.622	248	97992	37.712	ng	98
68) Hexachlorobenzene	10.686	284	114699	38.620	ng	97
69) Atrazine	10.792	200	82780	40.140	ng	97
70) Pentachlorophenol	10.916	266	13732m	28.402	ng	
71) Phenanthrene	11.104	178	482600	40.740	ng	99
72) Anthracene	11.157	178	470722	39.821	ng	99
73) Carbazole	11.322	167	426852	39.110	ng	98
74) Di-n-butylphthalate	11.639	149	504649	37.593	ng	99
75) Fluoranthene	12.286	202	489559	40.880	ng	100
77) Benzidine	12.422	184	106623	47.946	ng	99
78) Pyrene	12.516	202	494797	42.335	ng	99
80) Butylbenzylphthalate	13.133	149	186569	40.229	ng	98
81) Benzo(a)anthracene	13.704	228	369667	39.732	ng	99
82) 3,3'-Dichlorobenzidine	13.669	252	107049	39.621	ng	98
83) Chrysene	13.739	228	355524	39.438	ng	100
84) Bis(2-ethylhexyl)phtha...	13.686	149	212236	38.632	ng	99
85) Di-n-octyl phthalate	14.298	149	271646	37.505	ng	100
87) Indeno(1,2,3-cd)pyrene	16.398	276	338385	41.979	ng	100
88) Benzo(b)fluoranthene	14.716	252	274056	36.622	ng	98
89) Benzo(k)fluoranthene	14.739	252	294018	39.106	ng	100
90) Benzo(a)pyrene	15.039	252	235830	39.236	ng	99
91) Dibenzo(a,h)anthracene	16.398	278	280756	42.151	ng	100
92) Benzo(g,h,i)perylene	16.786	276	274605	41.107	ng	98

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

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