

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101422\
 Data File : BF130662.D
 Acq On : 14 Oct 2022 12:04
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 10/17/2022
 Supervised By : mohammad ahmed 10/18/2022

Quant Time: Oct 17 02:18:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 12 06:52:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.581	152	80182	20.000	ng	0.00
21) Naphthalene-d8	7.869	136	297979	20.000	ng	0.00
39) Acenaphthene-d10	9.616	164	159565	20.000	ng	0.00
64) Phenanthrene-d10	11.092	188	276517	20.000	ng	0.00
76) Chrysene-d12	13.727	240	169204	20.000	ng	0.00
86) Perylene-d12	15.098	264	128557	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.181	112	357813	80.439	ng	0.00
7) Phenol-d6	6.234	99	442803	79.175	ng	0.00
23) Nitrobenzene-d5	7.157	82	430218	78.817	ng	0.00
42) 2,4,6-Tribromophenol	10.404	330	132647	84.957	ng	0.00
45) 2-Fluorobiphenyl	8.945	172	860994	81.734	ng	0.00
79) Terphenyl-d14	12.680	244	849727	83.350	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.146	88	84916	31.553	ng	98
3) Pyridine	2.805	79	190125	39.495	ng	98
4) n-Nitrosodimethylamine	2.769	42	91802	40.475	ng	89
6) Aniline	6.251	93	261928	39.119	ng	# 83
8) 2-Chlorophenol	6.369	128	207806	39.725	ng	99
9) Benzaldehyde	6.134	77	126396	35.972	ng	99
10) Phenol	6.245	94	253809	39.222	ng	91
11) bis(2-Chloroethyl)ether	6.328	93	179797	38.250	ng	99
12) 1,3-Dichlorobenzene	6.522	146	228367	39.871	ng	99
13) 1,4-Dichlorobenzene	6.598	146	233470	40.180	ng	99
14) 1,2-Dichlorobenzene	6.751	146	217225	40.141	ng	99
15) Benzyl Alcohol	6.740	79	164399	40.943	ng	99
16) 2,2'-oxybis(1-Chloropr...	6.869	45	223629	36.979	ng	92
17) 2-Methylphenol	6.857	107	166168	39.933	ng	98
18) Hexachloroethane	7.087	117	85999	39.122	ng	95
19) n-Nitroso-di-n-propyla...	7.010	70	138228	38.308	ng	100
20) 3+4-Methylphenols	7.010	107	217025	38.922	ng	94
22) Acetophenone	7.004	105	283342	39.639	ng	99
24) Nitrobenzene	7.175	77	217836	39.684	ng	99
25) Isophorone	7.416	82	353081	38.865	ng	99
26) 2-Nitrophenol	7.492	139	110962	41.220	ng	98
27) 2,4-Dimethylphenol	7.534	122	151773	40.452	ng	99
28) bis(2-Chloroethoxy)met...	7.628	93	206081	39.177	ng	99
29) 2,4-Dichlorophenol	7.734	162	176441	42.032	ng	98
30) 1,2,4-Trichlorobenzene	7.810	180	187795	41.497	ng	98
31) Naphthalene	7.887	128	617721	40.047	ng	100
32) Benzoic acid	7.692	122	88069	37.103	ng	99
33) 4-Chloroaniline	7.945	127	242059	40.511	ng	99
34) Hexachlorobutadiene	8.004	225	120583	42.930	ng	97
35) Caprolactam	8.334	113	51192m	40.532	ng	
36) 4-Chloro-3-methylphenol	8.439	107	184614	40.544	ng	95
37) 2-Methylnaphthalene	8.581	142	403809	41.601	ng	99
38) 1-Methylnaphthalene	8.675	142	393724	41.783	ng	99
40) 1,2,4,5-Tetrachloroben...	8.745	216	177490	40.998	ng	98
41) Hexachlorocyclopentadiene	8.728	237	51444	37.180	ng	98
43) 2,4,6-Trichlorophenol	8.863	196	120580	41.482	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	8.910	196	131558	41.155	ng	95
46) 1,1'-Biphenyl	9.045	154	469192	40.574	ng	100
47) 2-Chloronaphthalene	9.063	162	386610	40.906	ng	97
48) 2-Nitroaniline	9.175	65	115177	40.483	ng	99
49) Acenaphthylene	9.475	152	575930	40.670	ng	99
50) Dimethylphthalate	9.351	163	440700	39.898	ng	99
51) 2,6-Dinitrotoluene	9.416	165	97968	40.434	ng	99
52) Acenaphthene	9.651	154	350683	40.874	ng	99
53) 3-Nitroaniline	9.586	138	107172	40.352	ng	96
54) 2,4-Dinitrophenol	9.698	184	40195	35.505	ng	# 88
55) Dibenzofuran	9.822	168	527894	40.389	ng	98
56) 4-Nitrophenol	9.763	139	56022	38.256	ng	93
57) 2,4-Dinitrotoluene	9.822	165	129480	41.581	ng	90
58) Fluorene	10.163	166	433651	40.357	ng	99
59) 2,3,4,6-Tetrachlorophenol	9.945	232	102905	40.379	ng	99
60) Diethylphthalate	10.045	149	429970	39.015	ng	99
61) 4-Chlorophenyl-phenyle...	10.157	204	196637	40.476	ng	98
62) 4-Nitroaniline	10.198	138	102187	40.299	ng	96
63) Azobenzene	10.316	77	395057	38.060	ng	98
65) 4,6-Dinitro-2-methylph...	10.228	198	72554	41.987	ng	89
66) n-Nitrosodiphenylamine	10.280	169	359341	39.067	ng	98
67) 4-Bromophenyl-phenylether	10.645	248	127684	40.810	ng	97
68) Hexachlorobenzene	10.704	284	147196	41.618	ng	97
69) Atrazine	10.810	200	104138	41.577	ng	99
70) Pentachlorophenol	10.910	266	47328	34.382	ng	97
71) Phenanthrene	11.122	178	605821	39.544	ng	99
72) Anthracene	11.169	178	589867	39.424	ng	100
73) Carbazole	11.333	167	530916	39.766	ng	99
74) Di-n-butylphthalate	11.663	149	639985	39.572	ng	100
75) Fluoranthene	12.304	202	610283	41.448	ng	99
77) Benzidine	12.433	184	145734	34.075	ng	99
78) Pyrene	12.527	202	605875	41.506	ng	100
80) Butylbenzylphthalate	13.157	149	225686	40.207	ng	100
81) Benzo(a)anthracene	13.716	228	459999	40.347	ng	99
82) 3,3'-Dichlorobenzidine	13.686	252	126188	38.301	ng	# 97
83) Chrysene	13.751	228	423981	39.222	ng	100
84) Bis(2-ethylhexyl)phtha...	13.716	149	261043	38.361	ng	99
85) Di-n-octyl phthalate	14.327	149	350389	33.621	ng	99
87) Indeno(1,2,3-cd)pyrene	16.392	276	412410	42.861	ng	98
88) Benzo(b)fluoranthene	14.721	252	334671	41.112	ng	98
89) Benzo(k)fluoranthene	14.745	252	324969	39.206	ng	99
90) Benzo(a)pyrene	15.045	252	270159	39.875	ng	99
91) Dibenzo(a,h)anthracene	16.398	278	338179	42.877	ng	99
92) Benzo(g,h,i)perylene	16.774	276	344902	46.951	ng	97

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

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