

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032919\
 Data File : BF113407.D
 Acq On : 29 Mar 2019 11:09
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 4/1/2019 5:46:23 PM

Quant Time: Mar 29 13:06:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	197252	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	788955	20.00	ng	0.00
39) Acenaphthene-d10	9.74	164	393678	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	789122	20.00	ng	0.00
76) Chrysene-d12	13.85	240	585092	20.00	ng	0.00
87) Perylene-d12	15.24	264	469804	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.32	112	869360	72.07	ng	0.00
7) Phenol-d6	6.36	99	1082393	70.93	ng	0.00
23) Nitrobenzene-d5	7.27	82	1018720	76.64	ng	0.00
42) 2,4,6-Tribromophenol	10.53	330	357229	73.46	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	1846003	70.48	ng	0.00
79) Terphenyl-d14	12.80	244	2029279	76.45	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.36	88	207844	33.958	ng	98
3) Pyridine	3.06	79	566166	37.522	ng	97
4) n-Nitrosodimethylamine	3.01	42	228322	34.038	ng	89
6) Aniline	6.37	93	671196	36.116	ng	97
8) 2-Chlorophenol	6.49	128	510323	36.429	ng	99
9) Benzaldehyde	6.25	77	371087	36.236	ng	99
10) Phenol	6.37	94	609245	34.969	ng	98
11) bis(2-Chloroethyl)ether	6.45	93	492335	36.865	ng	99
12) 1,3-Dichlorobenzene	6.65	146	545842	36.147	ng	98
13) 1,4-Dichlorobenzene	6.72	146	550913	37.785	ng	100
14) 1,2-Dichlorobenzene	6.87	146	500551	34.896	ng	99
15) Benzyl Alcohol	6.86	79	373974	37.145	ng	99
16) 2,2'-oxybis(1-Chloropropan	6.99	45	766949	38.228	ng	97
17) 2-Methylphenol	6.97	107	406488	37.026	ng	99
18) Hexachloroethane	7.22	117	200986	38.380	ng	95
19) n-Nitroso-di-n-propylamine	7.13	70	339522	35.419	ng	98
20) 3+4-Methylphenols	7.13	107	480695	36.746	ng	96
22) Acetophenone	7.12	105	658147	36.579	ng	98
24) Nitrobenzene	7.29	77	529035	38.140	ng	100
25) Isophorone	7.53	82	989049	38.395	ng	99
26) 2-Nitrophenol	7.60	139	291440	39.750	ng	97
27) 2,4-Dimethylphenol	7.65	122	410517	38.924	ng	100
28) bis(2-Chloroethoxy)methane	7.74	93	635289	39.162	ng	99
29) 2,4-Dichlorophenol	7.86	162	445122	38.946	ng	100
30) 1,2,4-Trichlorobenzene	7.93	180	469155	38.047	ng	98
31) Naphthalene	8.01	128	1443313	37.436	ng	100
32) Benzoic acid	7.80	122	309837m	36.496	ng	
33) 4-Chloroaniline	8.06	127	617516	41.074	ng	98
34) Hexachlorobutadiene	8.13	225	286143	39.043	ng	98
35) Caprolactam	8.45	113	148442	40.468	ng	96
36) 4-Chloro-3-methylphenol	8.56	107	452740	39.384	ng	98
37) 2-Methylnaphthalene	8.70	142	955788	39.731	ng	99
38) 1-Methylnaphthalene	8.80	142	923904	39.928	ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.87	216	464855	36.785	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.86	237	219427	41.023	ng	99
43) 2,4,6-Trichlorophenol	8.99	196	300917	36.567	ng	100
44) 2,4,5-Trichlorophenol	9.03	196	333701	36.642	ng	99
46) 1,1'-Biphenyl	9.16	154	1190150	36.197	ng	99
47) 2-Chloronaphthalene	9.19	162	917802	36.254	ng	99
48) 2-Nitroaniline	9.29	65	291570	36.701	ng	98
49) Acenaphthylene	9.60	152	1499567	36.634	ng	99
50) Dimethylphthalate	9.47	163	1102822	37.801	ng	100
51) 2,6-Dinitrotoluene	9.53	165	249863	37.963	ng #	84
52) Acenaphthene	9.77	154	837626	36.337	ng	99
53) 3-Nitroaniline	9.70	138	262840	35.412	ng	92
54) 2,4-Dinitrophenol	9.81	184	135875	41.507	ng	96
55) Dibenzofuran	9.95	168	1253080	36.159	ng	99
56) 4-Nitrophenol	9.87	139	172139	32.530	ng	94
57) 2,4-Dinitrotoluene	9.93	165	306061	36.567	ng	95
58) Fluorene	10.29	166	966838	35.914	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.07	232	297807	38.669	ng	97
60) Diethylphthalate	10.16	149	1061682	37.096	ng	99
61) 4-Chlorophenyl-phenylether	10.27	204	489649	37.318	ng	97
62) 4-Nitroaniline	10.31	138	290510	38.015	ng	98
63) Azobenzene	10.43	77	1021519	37.917	ng	97
65) 4,6-Dinitro-2-methylphenol	10.35	198	148869	34.486	ng	95
66) n-Nitrosodiphenylamine	10.40	169	883085	36.469	ng	100
67) 4-Bromophenyl-phenylether	10.76	248	330512	38.428	ng	97
68) Hexachlorobenzene	10.84	284	379262	37.990	ng	99
69) Atrazine	10.93	200	297869	38.973	ng	97
70) Pentachlorophenol	11.03	266	195731	37.631	ng	99
71) Phenanthrene	11.24	178	1433200	36.572	ng	100
72) Anthracene	11.29	178	1468867	36.902	ng	99
73) Carbazole	11.45	167	1428600	37.612	ng	99
74) Di-n-butylphthalate	11.78	149	1698829	38.570	ng	100
75) Fluoranthene	12.43	202	1573620	35.521	ng	99
77) Benzidine	12.55	184	525898	23.471	ng	99
78) Pyrene	12.65	202	1585911	39.115	ng	99
80) Butylbenzylphthalate	13.27	149	707746	39.608	ng	98
81) Benzo(a)anthracene	13.84	228	1216682	36.328	ng	100
82) 3,3'-Dichlorobenzidine	13.80	252	531586	39.565	ng	99
83) Chrysene	13.87	228	1280363	37.637	ng	99
84) Bis(2-ethylhexyl)phthalate	13.83	149	866970	39.579	ng	98
85) Di-n-octyl phthalate	14.43	149	1466989	39.454	ng	100
86) Indeno(1,2,3-cd)pyrene	16.61	276	1004171	34.395	ng	98
88) Benzo(b)fluoranthene	14.85	252	1076559	36.889	ng	99
89) Benzo(k)fluoranthene	14.88	252	1012167	39.493	ng	99
90) Benzo(a)pyrene	15.19	252	948919	36.676	ng	99
91) Dibenzo(a,h)anthracene	16.62	278	835672	37.355	ng	100
92) Benzo(g,h,i)perylene	17.01	276	836101	37.067	ng	98

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

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