

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010421\
 Data File : BF122909.D
 Acq On : 4 Jan 2021 14:45
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad

Quant Time: Jan 04 15:35:29 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF120520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 31 15:53:00 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.822	152	129680	20.00	ng	0.00
21) Naphthalene-d8	8.110	136	473650	20.00	ng	0.00
39) Acenaphthene-d10	9.868	164	249000	20.00	ng	0.00
64) Phenanthrene-d10	11.362	188	451101	20.00	ng	0.00
76) Chrysene-d12	14.015	240	344248	20.00	ng	0.00
86) Perylene-d12	15.486	264	321534	20.00	ng	0.00

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System Monitoring Compounds						
5) 2-Fluorophenol	5.439	112	639370	78.06	ng	0.00
7) Phenol-d6	6.469	99	806772	74.26	ng	0.00
23) Nitrobenzene-d5	7.392	82	698955	73.93	ng	0.00
42) 2,4,6-Tribromophenol	10.668	330	239243	73.40	ng	0.00
45) 2-Fluorobiphenyl	9.192	172	1273248	69.16	ng	0.00
79) Terphenyl-d14	12.951	244	1467792	74.63	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	2.563	88	145649	37.83	ng	97
3) Pyridine	3.304	79	378049	37.15	ng	98
4) n-Nitrosodimethylamine	3.263	42	136508	33.45	ng	99
6) Aniline	6.486	93	460405	36.00	ng	# 74
8) 2-Chlorophenol	6.616	128	346055	38.33	ng	95
9) Benzaldehyde	6.375	77	253279	38.52	ng	98
10) Phenol	6.480	94	411412	36.55	ng	88
11) bis(2-Chloroethyl)ether	6.563	93	319479	37.15	ng	98
12) 1,3-Dichlorobenzene	6.763	146	406761	38.08	ng	99
13) 1,4-Dichlorobenzene	6.839	146	410744	38.13	ng	97
14) 1,2-Dichlorobenzene	6.998	146	372553	35.83	ng	98
15) Benzyl Alcohol	6.975	79	274011	36.63	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.104	45	399397	32.57	ng	90
17) 2-Methylphenol	7.086	107	274126	38.20	ng	97
18) Hexachloroethane	7.333	117	144624	37.68	ng	97
19) n-Nitroso-di-n-propyla...	7.245	70	214570	34.48	ng	99
20) 3+4-Methylphenols	7.245	107	335737	37.03	ng	# 86
22) Acetophenone	7.239	105	438586	37.24	ng	99
24) Nitrobenzene	7.410	77	346992	36.90	ng	99
25) Isophorone	7.651	82	582164	35.75	ng	100
26) 2-Nitrophenol	7.727	139	185828	40.51	ng	95
27) 2,4-Dimethylphenol	7.769	122	251026	37.34	ng	99
28) bis(2-Chloroethoxy)met...	7.857	93	407571	36.55	ng	99
29) 2,4-Dichlorophenol	7.974	162	295034	37.58	ng	99
30) 1,2,4-Trichlorobenzene	8.057	180	328938	36.98	ng	97
31) Naphthalene	8.133	128	977801	37.41	ng	100
32) Benzoic acid	7.916	122	168815	31.51	ng	93
33) 4-Chloroaniline	8.186	127	410637	37.58	ng	99
34) Hexachlorobutadiene	8.251	225	202327	36.96	ng	98
35) Caprolactam	8.569	113	89289m	37.76	ng	
36) 4-Chloro-3-methylphenol	8.674	107	290544	37.57	ng	100
37) 2-Methylnaphthalene	8.827	142	640806	37.06	ng	99
38) 1-Methylnaphthalene	8.927	142	601133	36.75	ng	100
40) 1,2,4,5-Tetrachloroben...	8.992	216	310246	37.62	ng	100
41) Hexachlorocyclopentadiene	8.980	237	140619	38.56	ng	100
43) 2,4,6-Trichlorophenol	9.110	196	207493	36.43	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.157	196	213290	36.23	ng	96
46) 1,1'-Biphenyl	9.292	154	767223	37.13	ng	99
47) 2-Chloronaphthalene	9.316	162	642201	37.51	ng	99
48) 2-Nitroaniline	9.416	65	186867	37.44	ng	92
49) Acenaphthylene	9.733	152	940437	37.19	ng	99
50) Dimethylphthalate	9.592	163	746412	37.53	ng	100
51) 2,6-Dinitrotoluene	9.663	165	164405	39.14	ng	93
52) Acenaphthene	9.904	154	566739	34.64	ng	98
53) 3-Nitroaniline	9.833	138	185869	38.30	ng	100
54) 2,4-Dinitrophenol	9.945	184	98777	48.10	ng	96
55) Dibenzofuran	10.080	168	833632	36.22	ng	98
56) 4-Nitrophenol	10.004	139	134028	38.73	ng	91
57) 2,4-Dinitrotoluene	10.068	165	210262	37.58	ng	98
58) Fluorene	10.421	166	665661	36.36	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.204	232	183145	35.40	ng	100
60) Diethylphthalate	10.292	149	708260	36.62	ng	99
61) 4-Chlorophenyl-phenyle...	10.410	204	335119	37.18	ng	99
62) 4-Nitroaniline	10.451	138	191199	38.81	ng	95
63) Azobenzene	10.574	77	632401	35.57	ng	94
65) 4,6-Dinitro-2-methylph...	10.486	198	107287	39.00	ng	90
66) n-Nitrosodiphenylamine	10.533	169	588339	36.91	ng	100
67) 4-Bromophenyl-phenylether	10.904	248	228371	37.36	ng	93
68) Hexachlorobenzene	10.974	284	250925	37.13	ng	99
69) Atrazine	11.062	200	168814	32.61	ng	98
70) Pentachlorophenol	11.174	266	135366	37.81	ng	98
71) Phenanthrene	11.386	178	986038	36.78	ng	100
72) Anthracene	11.439	178	996727	37.49	ng	99
73) Carbazole	11.598	167	903274	37.47	ng	98
74) Di-n-butylphthalate	11.921	149	1096829	36.21	ng	99
75) Fluoranthene	12.580	202	1021868	36.35	ng	100
77) Benzidine	12.704	184	340839	45.77	ng	99
78) Pyrene	12.809	202	1017815	38.24	ng	99
80) Butylbenzylphthalate	13.427	149	444567	37.08	ng	92
81) Benzo(a)anthracene	14.003	228	877189	38.13	ng	98
82) 3,3'-Dichlorobenzidine	13.962	252	302413	36.70	ng	100
83) Chrysene	14.045	228	857368	37.45	ng	99
84) Bis(2-ethylhexyl)phtha...	13.986	149	589181	35.89	ng	99
85) Di-n-octyl phthalate	14.598	149	922846	33.89	ng	98
87) Indeno(1,2,3-cd)pyrene	16.974	276	832316	39.44	ng	100
88) Benzo(b)fluoranthene	15.062	252	808776	35.85	ng	98
89) Benzo(k)fluoranthene	15.092	252	795110	38.55	ng	99
90) Benzo(a)pyrene	15.427	252	736684	37.33	ng	99
91) Dibenzo(a,h)anthracene	16.980	278	683895	39.59	ng	99
92) Benzo(g,h,i)perylene	17.415	276	681477	40.76	ng	99

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

