

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041320\
 Data File : BF119968.D
 Acq On : 14 Apr 2020 22:35
 Operator : MA/SJ
 Sample : L2229-16MS
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-32MS

Manual Integrations
 APPROVED

mohammad
 4/16/2020 12:51:29 PM

Quant Time: Apr 15 04:07:38 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 14 12:44:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	200398	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	781446	20.00	ng	0.00
39) Acenaphthene-d10	9.78	164	408779	20.00	ng	0.00
64) Phenanthrene-d10	11.26	188	808915	20.00	ng	0.00
76) Chrysene-d12	13.91	240	560910	20.00	ng	0.00
87) Perylene-d12	15.33	264	468134	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	753493	64.98	ng	0.01
7) Phenol-d6	6.37	99	961964	64.38	ng	0.00
23) Nitrobenzene-d5	7.30	82	983900	65.14	ng	0.00
42) 2,4,6-Tribromophenol	10.57	330	379602	75.85	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	2080021	72.24	ng	0.00
79) Terphenyl-d14	12.86	244	2610946	78.32	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.52	88	139880	27.083	ng	95
3) Pyridine	3.19	79	323827	22.852	ng	95
4) n-Nitrosodimethylamine	3.13	42	212317	29.325	ng	94
6) Aniline	6.39	93	374912	19.117	ng	92
8) 2-Chlorophenol	6.52	128	321924	24.847	ng	96
9) Benzaldehyde	6.28	77	167768	15.168	ng	98
10) Phenol	6.38	94	383682	22.634	ng	90
11) bis(2-Chloroethyl)ether	6.47	93	454190	34.701	ng	97
12) 1,3-Dichlorobenzene	6.67	146	422681	29.799	ng	98
13) 1,4-Dichlorobenzene	6.75	146	436572	30.214	ng	99
14) 1,2-Dichlorobenzene	6.90	146	411344	30.890	ng	98
15) Benzyl Alcohol	6.88	79	385459	33.214	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.02	45	788445	30.957	ng	96
17) 2-Methylphenol	7.00	107	305670	26.436	ng	97
18) Hexachloroethane	7.25	117	157950	29.250	ng	92
19) n-Nitroso-di-n-propylamine	7.16	70	351319	36.564	ng	96
20) 3+4-Methylphenols	7.15	107	363070	25.674	ng	# 74
22) Acetophenone	7.15	105	659626	36.047	ng	# 96
24) Nitrobenzene	7.32	77	519443	34.185	ng	98
25) Isophorone	7.56	82	1011796	34.748	ng	100
26) 2-Nitrophenol	7.64	139	178637	23.531	ng	89
27) 2,4-Dimethylphenol	7.68	122	364740	31.856	ng	98
28) bis(2-Chloroethoxy)methane	7.77	93	629955	36.765	ng	98
29) 2,4-Dichlorophenol	7.88	162	305631	26.042	ng	98
30) 1,2,4-Trichlorobenzene	7.96	180	425260	31.980	ng	99
31) Naphthalene	8.05	128	1379298	33.548	ng	100
32) Benzoic acid	7.79	122	187915	18.688	ng	97
33) 4-Chloroaniline	8.09	127	332738	18.590	ng	97
34) Hexachlorobutadiene	8.16	225	233293	29.307	ng	99
35) Caprolactam	8.46	113	119307m	33.233	ng	
36) 4-Chloro-3-methylphenol	8.58	107	380303	29.931	ng	97
37) 2-Methylnaphthalene	8.73	142	1010232	36.243	ng	98
38) 1-Methylnaphthalene	8.83	142	967395	36.821	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.90	216	447609	34.669	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.89	237	399448	54.408	ng	97
43) 2,4,6-Trichlorophenol	9.02	196	237939	26.688	ng	98
44) 2,4,5-Trichlorophenol	9.06	196	268208	26.710	ng	97
46) 1,1'-Biphenyl	9.20	154	1300524	37.693	ng	99
47) 2-Chloronaphthalene	9.23	162	992228	37.331	ng	98
48) 2-Nitroaniline	9.32	65	393571	40.861	ng	96
49) Acenaphthylene	9.65	152	1631803	40.369	ng	99
50) Dimethylphthalate	9.51	163	1431877	45.286	ng	99
51) 2,6-Dinitrotoluene	9.57	165	294081	42.473	ng	93
52) Acenaphthene	9.82	154	1064044m	39.461	ng	
53) 3-Nitroaniline	9.73	138	234834	29.697	ng	96
54) 2,4-Dinitrophenol	9.85	184	150092	36.207	ng	95
55) Dibenzofuran	9.99	168	1506953	40.726	ng	100
56) 4-Nitrophenol	9.90	139	301715	51.279	ng	91
57) 2,4-Dinitrotoluene	9.97	165	390741	42.833	ng	99
58) Fluorene	10.33	166	1183451	39.992	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.10	232	238137	31.007	ng	97
60) Diethylphthalate	10.21	149	1339973	40.890	ng	99
61) 4-Chlorophenyl-phenylether	10.32	204	566344	37.340	ng	96
62) 4-Nitroaniline	10.36	138	316074	41.941	ng	96
63) Azobenzene	10.49	77	1240717	42.247	ng	96
65) 4,6-Dinitro-2-methylphenol	10.38	198	123036	23.087	ng	90
66) n-Nitrosodiphenylamine	10.45	169	1111144	41.303	ng	99
67) 4-Bromophenyl-phenylether	10.81	248	365501	36.333	ng	95
68) Hexachlorobenzene	10.88	284	392915	38.502	ng	94
69) Atrazine	10.97	200	348628	46.577	ng	97
70) Pentachlorophenol	11.07	266	287054	48.810	ng	98
71) Phenanthrene	11.29	178	1765515	41.010	ng	98
72) Anthracene	11.34	178	1833060	41.680	ng	100
73) Carbazole	11.50	167	1710197	41.120	ng	99
74) Di-n-butylphthalate	11.83	149	2150664	39.336	ng	99
75) Fluoranthene	12.48	202	1933772	41.630	ng	98
77) Benzidine	12.60	184	202630	10.687	ng	98
78) Pyrene	12.71	202	1958598	41.420	ng	99
80) Butylbenzylphthalate	13.33	149	909764	39.650	ng	96
81) Benzo(a)anthracene	13.90	228	1521734	41.239	ng	99
82) 3,3'-Dichlorobenzidine	13.86	252	398488	28.942	ng	98
83) Chrysene	13.93	228	1582228	42.200	ng	98
84) Bis(2-ethylhexyl)phthalate	13.89	149	1197363	38.535	ng	99
85) Di-n-octyl phthalate	14.50	149	2142969	40.506	ng	99
86) Indeno(1,2,3-cd)pyrene	16.72	276	1304650	39.474	ng	97
88) Benzo(b)fluoranthene	14.93	252	1488429	44.198	ng	99
89) Benzo(k)fluoranthene	14.96	252	1285772	44.251	ng	99
90) Benzo(a)pyrene	15.27	252	1219535	43.070	ng	99
91) Dibenzo(a,h)anthracene	16.74	278	1095432	43.675	ng	99
92) Benzo(g,h,i)perylene	17.14	276	1059210	42.783	ng	# 95

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

