

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081619\
 Data File : BF116309.D
 Acq On : 16 Aug 2019 9:17
 Operator : HP/JU
 Sample : PB122249BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB122249BS

Manual Integrations
 APPROVED

mohammad
 8/20/2019 8:36:43 AM

Quant Time: Aug 16 11:53:23 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 11:13:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	124478	20.00	ng	-0.03
21) Naphthalene-d8	8.10	136	515283	20.00	ng	-0.03
39) Acenaphthene-d10	9.85	164	274070	20.00	ng	-0.04
64) Phenanthrene-d10	11.34	188	469657	20.00	ng	-0.03
76) Chrysene-d12	13.98	240	329723	20.00	ng	-0.03
87) Perylene-d12	15.44	264	272322	20.00	ng	-0.04

System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	893640	120.18	ng	-0.01
7) Phenol-d6	6.46	99	1081954	115.33	ng	-0.02
23) Nitrobenzene-d5	7.39	82	714688	80.06	ng	-0.03
42) 2,4,6-Tribromophenol	10.64	330	301320	113.40	ng	-0.03
45) 2-Fluorobiphenyl	9.17	172	1215060	83.04	ng	-0.03
79) Terphenyl-d14	12.92	244	1346036	86.02	ng	-0.04

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.66	88	156197	41.581	ng	97
3) Pyridine	3.41	79	360398	34.345	ng	99
4) n-Nitrosodimethylamine	3.36	42	164632	39.756	ng	99
6) Aniline	6.49	93	401304	29.869	ng	# 86
8) 2-Chlorophenol	6.61	128	377869	45.956	ng	98
9) Benzaldehyde	6.38	77	127223	19.654	ng	98
10) Phenol	6.48	94	474828	42.980	ng	98
11) bis(2-Chloroethyl)ether	6.56	93	360961	44.441	ng	100
12) 1,3-Dichlorobenzene	6.76	146	401445	42.246	ng	100
13) 1,4-Dichlorobenzene	6.83	146	411396	43.238	ng	99
14) 1,2-Dichlorobenzene	6.99	146	376323	42.955	ng	98
15) Benzyl Alcohol	6.97	79	297977	41.853	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.09	45	473548	42.744	ng	83
17) 2-Methylphenol	7.08	107	310757	44.634	ng	97
18) Hexachloroethane	7.32	117	148464	42.381	ng	97
19) n-Nitroso-di-n-propylamine	7.23	70	255191	43.897	ng	97
20) 3+4-Methylphenols	7.23	107	388266	47.125	ng	92
22) Acetophenone	7.23	105	508283	44.977	ng	96
24) Nitrobenzene	7.40	77	420481	43.623	ng	99
25) Isophorone	7.64	82	755211	44.244	ng	98
26) 2-Nitrophenol	7.72	139	205168	45.156	ng	97
27) 2,4-Dimethylphenol	7.76	122	330566	48.358	ng	99
28) bis(2-Chloroethoxy)methane	7.85	93	464562	43.910	ng	100
29) 2,4-Dichlorophenol	7.97	162	326272	45.205	ng	99
30) 1,2,4-Trichlorobenzene	8.04	180	348687	42.381	ng	99
31) Naphthalene	8.12	128	1075638	44.360	ng	99
32) Benzoic acid	7.92	122	193804	40.213	ng	96
33) 4-Chloroaniline	8.17	127	300203	27.709	ng	99
34) Hexachlorobutadiene	8.23	225	196574	41.480	ng	99
35) Caprolactam	8.55	113	98086m	42.018	ng	
36) 4-Chloro-3-methylphenol	8.66	107	335694	44.708	ng	99
37) 2-Methylnaphthalene	8.81	142	716890	44.891	ng	99
38) 1-Methylnaphthalene	8.91	142	666498	44.116	ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.98	216	330016	45.041	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.96	237	135192	43.980	ng	99
43) 2,4,6-Trichlorophenol	9.10	196	207799	41.203	ng	98
44) 2,4,5-Trichlorophenol	9.15	196	219906	42.182	ng	99
46) 1,1'-Biphenyl	9.27	154	874448	45.193	ng	97
47) 2-Chloronaphthalene	9.30	162	683848	42.464	ng	99
48) 2-Nitroaniline	9.40	65	216100	40.597	ng	96
49) Acenaphthylene	9.72	152	1008196	42.912	ng	99
50) Dimethylphthalate	9.57	163	788600	42.259	ng	100
51) 2,6-Dinitrotoluene	9.65	165	175498	43.275	ng	97
52) Acenaphthene	9.89	154	620764	43.068	ng	99
53) 3-Nitroaniline	9.82	138	150709	30.076	ng	100
54) 2,4-Dinitrophenol	9.94	184	124106	62.106	ng	93
55) Dibenzofuran	10.06	168	876526	41.817	ng	98
56) 4-Nitrophenol	10.00	139	214597	66.852	ng	91
57) 2,4-Dinitrotoluene	10.06	165	217501	40.282	ng	93
58) Fluorene	10.40	166	721356	44.730	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.19	232	187175	42.676	ng	98
60) Diethylphthalate	10.27	149	760070	43.626	ng	99
61) 4-Chlorophenyl-phenylether	10.39	204	364604	44.641	ng	98
62) 4-Nitroaniline	10.44	138	173759	33.554	ng	96
63) Azobenzene	10.55	77	791974	44.195	ng	98
65) 4,6-Dinitro-2-methylphenol	10.47	198	108351	43.534	ng	90
66) n-Nitrosodiphenylamine	10.51	169	640827	45.332	ng	99
67) 4-Bromophenyl-phenylether	10.87	248	220988	43.954	ng	95
68) Hexachlorobenzene	10.96	284	246584	42.414	ng	95
69) Atrazine	11.04	200	196457	42.244	ng	98
70) Pentachlorophenol	11.16	266	188090	66.639	ng	97
71) Phenanthrene	11.36	178	1045708	43.553	ng	100
72) Anthracene	11.42	178	1090785	44.890	ng	99
73) Carbazole	11.57	167	975270	44.239	ng	99
74) Di-n-butylphthalate	11.89	149	1200214	46.670	ng	98
75) Fluoranthene	12.55	202	1112666	44.266	ng	98
77) Benzidine	12.67	184	289361	25.976	ng	97
78) Pyrene	12.78	202	1113085	44.783	ng	99
80) Butylbenzylphthalate	13.39	149	504127	47.949	ng	99
81) Benzo(a)anthracene	13.97	228	906987	43.037	ng	99
82) 3,3'-Dichlorobenzidine	13.93	252	292565	39.958	ng	99
83) Chrysene	14.01	228	902389	44.104	ng	99
84) Bis(2-ethylhexyl)phthalate	13.94	149	703016	51.456	ng	100
85) Di-n-octyl phthalate	14.55	149	1105403	50.938	ng	99
86) Indeno(1,2,3-cd)pyrene	16.90	276	765737	44.560	ng	99
88) Benzo(b)fluoranthene	15.02	252	848901	53.912	ng	99
89) Benzo(k)fluoranthene	15.04	252	705384	45.993	ng	98
90) Benzo(a)pyrene	15.38	252	760936	54.060	ng	99
91) Dibenzo(a,h)anthracene	16.90	278	661372	54.282	ng	99
92) Benzo(g,h,i)perylene	17.34	276	649362	54.268	ng	97

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

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