

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122220\
 Data File : BF122814.D
 Acq On : 22 Dec 2020 14:34
 Operator : JU/CG
 Sample : L5155-01MSD 2X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-04-12-18-20MSD

Manual Integrations
 APPROVED

mohammad
 12/23/2020 3:28:07 PM

Quant Time: Dec 22 15:08:11 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF120520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 15 18:14:23 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.822	152	136913	20.00	ng	0.00
21) Naphthalene-d8	8.104	136	504253	20.00	ng	0.00
39) Acenaphthene-d10	9.863	164	241681	20.00	ng	0.00
64) Phenanthrene-d10	11.345	188	395170	20.00	ng	0.00
76) Chrysene-d12	13.998	240	335696	20.00	ng	0.01
86) Perylene-d12	15.462	264	287151	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	363172	41.99	ng	0.00
7) Phenol-d6	6.451	99	429042	37.41	ng	0.00
23) Nitrobenzene-d5	7.381	82	262628	26.09	ng	-0.01
42) 2,4,6-Tribromophenol	10.651	330	109885	34.74	ng	0.00
45) 2-Fluorobiphenyl	9.181	172	481305	26.94	ng	0.00
79) Terphenyl-d14	12.945	244	459593	23.96	ng	0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.540	88	51354	12.63	ng	99
3) Pyridine	3.293	79	133746	12.45	ng	96
4) n-Nitrosodimethylamine	3.216	42	56189	13.04	ng	# 99
6) Aniline	6.481	93	78988	5.85	ng	# 91
8) 2-Chlorophenol	6.604	128	142725	14.97	ng	97
9) Benzaldehyde	6.369	77	41017	5.91	ng	94
10) Phenol	6.463	94	179009	15.06	ng	87
11) bis(2-Chloroethyl)ether	6.551	93	135856	14.96	ng	97
12) 1,3-Dichlorobenzene	6.757	146	159187	14.11	ng	99
13) 1,4-Dichlorobenzene	6.834	146	163631	14.39	ng	98
14) 1,2-Dichlorobenzene	6.987	146	152400	13.88	ng	99
15) Benzyl Alcohol	6.963	79	109441	13.86	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.098	45	166522	12.86	ng	95
17) 2-Methylphenol	7.075	107	111795	14.75	ng	98
18) Hexachloroethane	7.334	117	45771	11.29	ng	94
19) n-Nitroso-di-n-propyla...	7.228	70	95968	14.61	ng	97
20) 3+4-Methylphenols	7.234	107	143324	14.97	ng	# 85
22) Acetophenone	7.228	105	197286	15.74	ng	96
24) Nitrobenzene	7.398	77	140562	14.04	ng	99
25) Isophorone	7.639	82	248760	14.35	ng	99
26) 2-Nitrophenol	7.722	139	59450	12.17	ng	92
27) 2,4-Dimethylphenol	7.763	122	118872	16.61	ng	98
28) bis(2-Chloroethoxy)met...	7.851	93	164183	13.83	ng	99
29) 2,4-Dichlorophenol	7.969	162	113536	13.58	ng	98
30) 1,2,4-Trichlorobenzene	8.045	180	127619	13.48	ng	99
31) Naphthalene	8.128	128	401635	14.43	ng	99
32) Benzoic acid	7.851	122	55513	9.73	ng	97
33) 4-Chloroaniline	8.181	127	64695	5.56	ng	99
34) Hexachlorobutadiene	8.245	225	77140	13.23	ng	99
35) Caprolactam	8.522	113	37585	14.93	ng	89
36) 4-Chloro-3-methylphenol	8.663	107	110915	13.47	ng	99
37) 2-Methylnaphthalene	8.816	142	270174	14.68	ng	99
38) 1-Methylnaphthalene	8.916	142	243740	14.00	ng	99
40) 1,2,4,5-Tetrachloroben...	8.986	216	124589	15.56	ng	99
43) 2,4,6-Trichlorophenol	9.098	196	77398	14.00	ng	99
44) 2,4,5-Trichlorophenol	9.145	196	78991	13.82	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	9.281	154	313115	15.61	ng	97
47) 2-Chloronaphthalene	9.304	162	235880	14.19	ng	99
48) 2-Nitroaniline	9.404	65	68196	14.08	ng	98
49) Acenaphthylene	9.722	152	351409	14.32	ng	98
50) Dimethylphthalate	9.575	163	300483	15.57	ng	100
51) 2,6-Dinitrotoluene	9.639	165	51621	12.66	ng	91
52) Acenaphthene	9.892	154	210384	13.25	ng	100
53) 3-Nitroaniline	9.810	138	45019	9.56	ng	91
55) Dibenzofuran	10.063	168	320671	14.35	ng	98
56) 4-Nitrophenol	9.992	139	70206	20.90	ng	95
57) 2,4-Dinitrotoluene	10.051	165	63531	11.70	ng	# 88
58) Fluorene	10.404	166	247988	13.96	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.186	232	61327	12.21	ng	# 95
60) Diethylphthalate	10.275	149	269518	14.36	ng	99
61) 4-Chlorophenyl-phenyle...	10.398	204	122751	14.03	ng	98
62) 4-Nitroaniline	10.422	138	57038	11.93	ng	92
63) Azobenzene	10.557	77	222317	12.88	ng	98
66) n-Nitrosodiphenylamine	10.516	169	217478	15.58	ng	98
67) 4-Bromophenyl-phenylether	10.886	248	76342	14.26	ng	99
68) Hexachlorobenzene	10.963	284	80702	13.63	ng	# 90
69) Atrazine	11.045	200	64269	14.17	ng	100
70) Pentachlorophenol	11.157	266	73774	23.52	ng	98
71) Phenanthrene	11.369	178	350285	14.92	ng	99
72) Anthracene	11.422	178	352836	15.15	ng	98
73) Carbazole	11.580	167	301185	14.26	ng	98
74) Di-n-butylphthalate	11.910	149	392739	14.80	ng	99
75) Fluoranthene	12.574	202	289203	11.74	ng	95
77) Benzidine	12.727	184	40607	5.59	ng	# 91
78) Pyrene	12.839	202	383549m	14.78	ng	
80) Butylbenzylphthalate	13.410	149	149421	12.78	ng	96
81) Benzo(a)anthracene	13.986	228	329232	14.68	ng	98
82) 3,3'-Dichlorobenzidine	13.945	252	74186	9.23	ng	# 98
83) Chrysene	14.021	228	328389	14.71	ng	98
84) Bis(2-ethylhexyl)phtha...	13.968	149	232232	14.51	ng	100
85) Di-n-octyl phthalate	14.580	149	374100	14.09	ng	# 98
87) Indeno(1,2,3-cd)pyrene	16.921	276	271089	14.38	ng	99
88) Benzo(b)fluoranthene	15.033	252	287967	14.29	ng	96
89) Benzo(k)fluoranthene	15.062	252	294383	15.98	ng	94
90) Benzo(a)pyrene	15.398	252	249020	14.13	ng	# 95
91) Dibenzo(a,h)anthracene	16.933	278	222951	14.45	ng	97
92) Benzo(g,h,i)perylene	17.362	276	219244	14.68	ng	99

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

