

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122719\
 Data File : BF118340.D
 Acq On : 27 Dec 2019 16:17
 Operator : JU
 Sample : K6439-01MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LIED-BF001-01MSD

Manual Integrations
 APPROVED

mohammad
 12/29/2019 5:27:41 PM

Quant Time: Dec 28 01:31:12 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 15:58:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	156945	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	623815	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	340161	20.00	ng	0.00
64) Phenanthrene-d10	11.40	188	651709	20.00	ng	0.00
76) Chrysene-d12	14.05	240	452276	20.00	ng	0.00
87) Perylene-d12	15.54	264	449463	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.48	112	1115874	121.37	ng	0.02
7) Phenol-d6	6.50	99	1427780	125.18	ng	0.01
23) Nitrobenzene-d5	7.42	82	851335	78.06	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	514098	122.38	ng	0.01
45) 2-Fluorobiphenyl	9.22	172	1785273	80.11	ng	0.00
79) Terphenyl-d14	12.99	244	2251886	82.60	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.64	88	130831	35.795	ng	99
3) Pyridine	3.40	79	318952	32.753	ng	99
4) n-Nitrosodimethylamine	3.35	42	159546	39.627	ng	100
6) Aniline	6.52	93	432207	29.842	ng	89
8) 2-Chlorophenol	6.65	128	439653	42.221	ng	99
9) Benzaldehyde	6.40	77	244813	36.610	ng	98
10) Phenol	6.51	94	533335	41.603	ng	90
11) bis(2-Chloroethyl)ether	6.59	93	378072	41.238	ng	98
12) 1,3-Dichlorobenzene	6.79	146	473190	39.456	ng	99
13) 1,4-Dichlorobenzene	6.87	146	476690	39.133	ng	99
14) 1,2-Dichlorobenzene	7.02	146	454273	39.500	ng	99
15) Benzyl Alcohol	7.00	79	361109	41.582	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.13	45	389956	39.860	ng	93
17) 2-Methylphenol	7.12	107	348640	41.743	ng	95
18) Hexachloroethane	7.36	117	175193	39.187	ng	98
19) n-Nitroso-di-n-propylamine	7.27	70	288214	41.019	ng	97
20) 3+4-Methylphenols	7.27	107	448301	41.939	ng	91
22) Acetophenone	7.26	105	585900	38.322	ng	# 94
24) Nitrobenzene	7.44	77	426743	39.059	ng	97
25) Isophorone	7.68	82	797318	41.946	ng	100
26) 2-Nitrophenol	7.76	139	248754	42.985	ng	99
27) 2,4-Dimethylphenol	7.79	122	387806	49.116	ng	100
28) bis(2-Chloroethoxy)methane	7.89	93	493483	39.697	ng	100
29) 2,4-Dichlorophenol	8.00	162	379397	41.747	ng	97
30) 1,2,4-Trichlorobenzene	8.08	180	411795	38.898	ng	98
31) Naphthalene	8.16	128	1281168	40.407	ng	100
32) Benzoic acid	7.94	122	230052	36.100	ng	94
33) 4-Chloroaniline	8.21	127	219095	16.119	ng	98
34) Hexachlorobutadiene	8.27	225	240588	38.178	ng	99
35) Caprolactam	8.59	113	120633m	42.517	ng	
36) 4-Chloro-3-methylphenol	8.71	107	402360	41.237	ng	97
37) 2-Methylnaphthalene	8.86	142	891997	41.091	ng	100
38) 1-Methylnaphthalene	8.96	142	831672	40.664	ng	98
40) 1,2,4,5-Tetrachlorobenzene	9.03	216	391091	38.266	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122719\
 Data File : BF118340.D
 Acq On : 27 Dec 2019 16:17
 Operator : JU
 Sample : K6439-01MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LIED-BF001-01MSD

Manual Integrations
 APPROVED

mohammad
 12/29/2019 5:27:41 PM

Quant Time: Dec 28 01:31:12 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 15:58:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.01	237	308305	72.579	nq	98
43) 2,4,6-Trichlorophenol	9.14	196	271755	39.801	nq	99
44) 2,4,5-Trichlorophenol	9.19	196	285085	40.588	nq	99
46) 1,1'-Biphenyl	9.32	154	1066358	39.635	nq	99
47) 2-Chloronaphthalene	9.35	162	832444	38.443	nq	97
48) 2-Nitroaniline	9.45	65	233080	38.108	nq	90
49) Acenaphthylene	9.77	152	1330642	41.434	nq	100
50) Dimethylphthalate	9.62	163	1147657	44.947	nq	99
51) 2,6-Dinitrotoluene	9.69	165	225457	40.956	nq	93
52) Acenaphthene	9.94	154	765056	39.110	nq	99
53) 3-Nitroaniline	9.86	138	136563	21.136	nq	98
54) 2,4-Dinitrophenol	9.98	184	190161	69.256	nq	94
55) Dibenzofuran	10.11	168	1164609	40.292	nq	96
56) 4-Nitrophenol	10.04	139	333943	83.723	nq	91
57) 2,4-Dinitrotoluene	10.10	165	304132	41.346	nq	92
58) Fluorene	10.46	166	937149	40.164	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.23	232	237948	39.870	nq	99
60) Diethylphthalate	10.33	149	1036725	41.041	nq	100
61) 4-Chlorophenyl-phenylether	10.45	204	470876	40.426	nq	95
62) 4-Nitroaniline	10.49	138	242429	36.118	nq	95
63) Azobenzene	10.60	77	894111	41.323	nq	97
65) 4,6-Dinitro-2-methylphenol	10.52	198	148458	42.227	nq	94
66) n-Nitrosodiphenylamine	10.57	169	840821	40.255	nq	98
67) 4-Bromophenyl-phenylether	10.93	248	297636	39.025	nq	93
68) Hexachlorobenzene	11.01	284	341074	37.992	nq	94
69) Atrazine	11.10	200	294040	45.613	nq	96
70) Pentachlorophenol	11.21	266	306085	78.726	nq	99
71) Phenanthrene	11.42	178	1403978	39.518	nq	100
72) Anthracene	11.47	178	1472761	41.415	nq	100
73) Carbazole	11.63	167	1267265	40.161	nq	98
74) Di-n-butylphthalate	11.95	149	1701362	42.523	nq	99
75) Fluoranthene	12.62	202	1459132	40.695	nq	96
77) Benzidine	12.74	184	858987	69.314	nq	98
78) Pyrene	12.85	202	1453611	39.358	nq	99
80) Butylbenzylphthalate	13.46	149	712655	42.855	nq	95
81) Benzo(a)anthracene	14.04	228	1259472	39.784	nq	99
82) 3,3'-Dichlorobenzidine	14.00	252	334465	31.160	nq	98
83) Chrysene	14.08	228	1199810	39.580	nq	100
84) Bis(2-ethylhexyl)phthalate	14.02	149	1047734	47.915	nq	99
85) Di-n-octyl phthalate	14.63	149	1655431	51.467	nq	99
86) Indeno(1,2,3-cd)pyrene	17.07	276	1211661	46.911	nq	100
88) Benzo(b)fluoranthene	15.10	252	1114227	37.346	nq	99
89) Benzo(k)fluoranthene	15.13	252	1140120	39.799	nq	99
90) Benzo(a)pyrene	15.48	252	998324	37.984	nq	98
91) Dibenzo(a,h)anthracene	17.07	278	1024248	44.908	nq	99
92) Benzo(g,h,i)perylene	17.52	276	1001199	45.443	nq	98

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

