

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072220\
 Data File : BF120967.D
 Acq On : 22 Jul 2020 16:32
 Operator : JU/CG
 Sample : L3184-17MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-03-072020MS

Manual Integrations
 APPROVED

Sohil
 7/24/2020 12:10:42 PM

Quant Time: Jul 22 17:22:36 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 16 14:20:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	169613	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	662950	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	349694	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	671983	20.00	ng	0.00
76) Chrysene-d12	13.97	240	517663	20.00	ng	0.00
86) Perylene-d12	15.42	264	516063	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.45	112	1078519	104.24	ng	0.02
7) Phenol-d6	6.45	99	1230683	92.08	ng	0.00
23) Nitrobenzene-d5	7.38	82	934846	81.59	ng	0.00
42) 2,4,6-Tribromophenol	10.65	330	443092	115.76	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	1632475	69.70	ng	0.00
79) Terphenyl-d14	12.92	244	1760734	73.94	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.71	88	186037	39.069	ng	# 85
3) Pyridine	3.42	79	355097	28.033	ng	100
4) n-Nitrosodimethylamine	3.35	42	223144	41.197	ng	94
6) Aniline	6.47	93	328757	18.845	ng	# 90
8) 2-Chlorophenol	6.60	128	521210	45.729	ng	100
9) Benzaldehyde	6.36	77	234308	34.983	ng	98
10) Phenol	6.46	94	517586	35.206	ng	98
11) bis(2-Chloroethyl)ether	6.55	93	501291	44.491	ng	100
12) 1,3-Dichlorobenzene	6.76	146	467092	37.793	ng	99
13) 1,4-Dichlorobenzene	6.83	146	473444	38.524	ng	99
14) 1,2-Dichlorobenzene	6.99	146	453555	39.011	ng	99
15) Benzyl Alcohol	6.96	79	428293	45.315	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.09	45	675471	41.593	ng	99
17) 2-Methylphenol	7.07	107	411457	40.729	ng	99
18) Hexachloroethane	7.33	117	172035	38.676	ng	98
19) n-Nitroso-di-n-propylamine	7.23	70	328622	43.242	ng	98
20) 3+4-Methylphenols	7.23	107	447921	34.855	ng	# 82
22) Acetophenone	7.23	105	653501	43.922	ng	95
24) Nitrobenzene	7.40	77	524232	42.453	ng	98
25) Isophorone	7.64	82	1000395	43.750	ng	98
26) 2-Nitrophenol	7.72	139	277622	47.337	ng	98
27) 2,4-Dimethylphenol	7.75	122	459600	48.267	ng	99
28) bis(2-Chloroethoxy)methane	7.85	93	626959	44.918	ng	99
29) 2,4-Dichlorophenol	7.96	162	433884	44.592	ng	99
30) 1,2,4-Trichlorobenzene	8.04	180	427555	39.606	ng	99
31) Naphthalene	8.12	128	1374016	40.405	ng	100
32) Benzoic acid	7.86	122	158210	22.134	ng	96
33) 4-Chloroaniline	8.17	127	245848	16.206	ng	99
34) Hexachlorobutadiene	8.24	225	236132	37.105	ng	100
35) Caprolactam	8.55	113	104823m	33.594	ng	
36) 4-Chloro-3-methylphenol	8.66	107	455768	45.672	ng	94
37) 2-Methylnaphthalene	8.81	142	961563	43.548	ng	99
38) 1-Methylnaphthalene	8.91	142	906291	43.338	ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.98	216	429229	42.128	ng	99

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41) Hexachlorocyclopentadiene	8.96	237	422716	75.069	nq	98
43) 2,4,6-Trichlorophenol	9.09	196	324590	45.247	nq	98
44) 2,4,5-Trichlorophenol	9.13	196	347331	42.684	nq	96
46) 1,1'-Biphenyl	9.27	154	1165825	43.705	nq	99
47) 2-Chloronaphthalene	9.30	162	907445	41.726	nq	99
48) 2-Nitroaniline	9.40	65	296703	45.144	nq	91
49) Acenaphthylene	9.72	152	1464955	43.360	nq	100
50) Dimethylphthalate	9.58	163	1333861	52.872	nq	98
51) 2,6-Dinitrotoluene	9.64	165	254810	47.012	nq	94
52) Acenaphthene	9.89	154	929070	43.253	nq	100
53) 3-Nitroaniline	9.80	138	151915	22.348	nq	96
54) 2,4-Dinitrophenol	9.91	184	112074	39.688	nq	90
55) Dibenzofuran	10.06	168	1308345	42.120	nq	99
56) 4-Nitrophenol	9.97	139	386443	74.838	nq	100
57) 2,4-Dinitrotoluene	10.05	165	343629	48.278	nq	97
58) Fluorene	10.40	166	941614	40.644	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.17	232	300067	48.432	nq	96
60) Diethylphthalate	10.28	149	1145871	44.905	nq	98
61) 4-Chlorophenyl-phenylether	10.39	204	466400	37.745	nq	99
62) 4-Nitroaniline	10.42	138	270145	40.797	nq	99
63) Azobenzene	10.56	77	1054959	43.426	nq	99
65) 4,6-Dinitro-2-methylphenol	10.45	198	113282	30.638	nq	87
66) n-Nitrosodiphenylamine	10.52	169	945403	44.723	nq	100
67) 4-Bromophenyl-phenylether	10.88	248	325387	43.431	nq	99
68) Hexachlorobenzene	10.95	284	360626	44.497	nq	# 89
69) Atrazine	11.04	200	272897	52.126	nq	100
70) Pentachlorophenol	11.14	266	390183	84.038	nq	99
71) Phenanthrene	11.36	178	1477189	42.742	nq	100
72) Anthracene	11.42	178	1545491	44.534	nq	99
73) Carbazole	11.57	167	1461714	42.442	nq	99
74) Di-n-butylphthalate	11.90	149	1794056	45.141	nq	100
75) Fluoranthene	12.55	202	1653468	43.163	nq	98
77) Benzidine	12.67	184	490597	36.046	nq	99
78) Pyrene	12.78	202	1694752	46.306	nq	99
80) Butylbenzylphthalate	13.40	149	800410	49.059	nq	98
81) Benzo(a)anthracene	13.97	228	1324551	42.254	nq	99
82) 3,3'-Dichlorobenzidine	13.93	252	409803	34.652	nq	99
83) Chrysene	14.00	228	1454391	44.149	nq	99
84) Bis(2-ethylhexyl)phthalate	13.96	149	941494	46.797	nq	100
85) Di-n-octyl phthalate	14.57	149	1920695	47.986	nq	99
87) Indeno(1,2,3-cd)pyrene	16.85	276	1490020	41.516	nq	99
88) Benzo(b)fluoranthene	15.00	252	1478604	47.400	nq	100
89) Benzo(k)fluoranthene	15.03	252	1415526	49.757	nq	100
90) Benzo(a)pyrene	15.36	252	1297707	45.597	nq	99
91) Dibenzo(a,h)anthracene	16.87	278	1248684	42.592	nq	99
92) Benzo(g,h,i)perylene	17.28	276	1210640	41.881	nq	99

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

