

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072518\
 Data File : BF107652.D
 Acq On : 25 Jul 2018 13:05
 Operator : JU/SJ
 Sample : J4101-02MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1MS

Manual Integrations
 APPROVED

Sohil
 7/26/2018 2:02:40 PM

Quant Time: Jul 25 14:22:46 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 20 16:29:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	245552	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	945349	20.00	ng	-0.01
38) Acenaphthene-d10	10.01	164	442779	20.00	ng	0.00
63) Phenanthrene-d10	11.51	188	877123	20.00	ng	0.00
75) Chrysene-d12	14.16	240	649497	20.00	ng	0.00
86) Perylene-d12	15.68	264	426615	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	1615252	105.76	ng	0.02
7) Phenol-d6	6.61	99	1883196	102.69	ng	0.00
23) Nitrobenzene-d5	7.54	82	1386993	99.02	ng	-0.01
41) 2,4,6-Tribromophenol	10.81	330	696068	138.24	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	2569695	106.36	ng	-0.01
78) Terphenyl-d14	13.09	244	2566338	101.16	ng	-0.01

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.09	88	194081	27.303	ng	# 68
3) Pyridine	3.78	79	375263	19.405	ng	# 82
4) n-Nitrosodimethylamine	3.72	42	270009	33.105	ng	97
6) Aniline	6.65	93	138863	5.871	ng	# 56
8) 2-Chlorophenol	6.76	128	657644	40.193	ng	94
9) Benzaldehyde	6.53	77	63590	3.282	ng	86
10) Phenol	6.62	94	662058	30.698	ng	92
11) bis(2-Chloroethyl)ether	6.71	93	620741	34.716	ng	92
12) 1,3-Dichlorobenzene	6.91	146	722040	38.921	ng	97
13) 1,4-Dichlorobenzene	6.98	146	768009	40.392	ng	99
14) 1,2-Dichlorobenzene	7.14	146	678588	39.769	ng	99
15) Benzyl Alcohol	7.11	79	462288	36.989	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.24	45	1082042	35.871	ng	79
17) 2-Methylphenol	7.22	107	558392	40.152	ng	# 89
18) Hexachloroethane	7.47	117	270873	40.616	ng	97
19) n-Nitroso-di-n-propylamine	7.38	70	460958	38.910	ng	100
20) 3+4-Methylphenols	7.38	107	700857	42.441	ng	# 50
22) Acetophenone	7.38	105	954147	42.977	ng	# 90
24) Nitrobenzene	7.55	77	709827	44.242	ng	98
25) Isophorone	7.79	82	1292666	43.220	ng	97
26) 2-Nitrophenol	7.87	139	388603	57.356	ng	97
27) 2,4-Dimethylphenol	7.90	122	622337	48.526	ng	98
28) bis(2-Chloroethoxy)methane	7.99	93	808875	40.468	ng	99
29) 2,4-Dichlorophenol	8.11	162	602672	45.976	ng	98
30) 1,2,4-Trichlorobenzene	8.19	180	628601	42.696	ng	98
31) Naphthalene	8.27	128	1966961	47.316	ng	98
32) Benzoic acid	8.04	122	372939	42.698	ng	99
33) 4-Chloroaniline	8.34	127	47393	2.608	ng	95
34) Hexachlorobutadiene	8.38	225	372400	42.568	ng	99
35) Caprolactam	8.71	113	171126m	40.136	ng	
36) 4-Chloro-3-methylphenol	8.81	107	651981	44.776	ng	99
37) 2-Methylnaphthalene	8.97	142	1388556	48.202	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.13	216	651525	44.116	ng	98
40) Hexachlorocyclopentadiene	9.11	237	720151	95.925	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.25	196	414033	43.799	nq	97
43) 2,4,5-Trichlorophenol	9.30	196	456338	46.189	nq	97
45) 1,1'-Biphenyl	9.43	154	1680215	43.561	nq	98
46) 2-Chloronaphthalene	9.46	162	1267764	43.000	nq	99
47) 2-Nitroaniline	9.56	65	410937	47.810	nq	91
48) Acenaphthylene	9.88	152	2017154	45.545	nq	97
49) Dimethylphthalate	9.73	163	1445728	43.323	nq	98
50) 2,6-Dinitrotoluene	9.80	165	325246	49.090	nq	93
51) Acenaphthene	10.05	154	1141739	41.145	nq	99
52) 3-Nitroaniline	9.98	138	35341	4.385	nq	97
53) 2,4-Dinitrophenol	10.08	184	384175	105.911	nq #	21
54) Dibenzofuran	10.22	168	1741679	42.633	nq	96
55) 4-Nitrophenol	10.15	139	434441	71.990	nq	94
56) 2,4-Dinitrotoluene	10.21	165	403752	50.485	nq #	95
57) Fluorene	10.57	166	1382079	47.421	nq	95
58) 2,3,4,6-Tetrachlorophenol	10.34	232	398041	48.409	nq	87
59) Diethylphthalate	10.42	149	1458208	41.840	nq	97
60) 4-Chlorophenyl-phenylether	10.55	204	686632	41.673	nq	94
61) 4-Nitroaniline	10.60	138	224723	33.185	nq	95
62) Azobenzene	10.71	77	1343288	41.104	nq	93
64) 4,6-Dinitro-2-methylphenol	10.62	198	239790	55.069	nq	83
65) n-Nitrosodiphenylamine	10.67	169	910577	30.567	nq	99
66) 4-Bromophenyl-phenylether	11.04	248	428350	41.486	nq	91
67) Hexachlorobenzene	11.11	284	475077	41.817	nq #	85
68) Atrazine	11.20	200	159354	17.520	nq	95
69) Pentachlorophenol	11.31	266	507044	71.453	nq	99
70) Phenanthrene	11.54	178	1883424	44.064	nq	99
71) Anthracene	11.58	178	2040531	47.415	nq	98
72) Carbazole	11.74	167	1319593	29.487	nq	99
73) Di-n-butylphthalate	12.05	149	2132386	46.893	nq	99
74) Fluoranthene	12.72	202	2124264	46.315	nq	97
77) Pyrene	12.95	202	2079259	46.802	nq	98
79) Butylbenzylphthalate	13.56	149	942457	49.328	nq #	91
80) Benzo(a)anthracene	14.15	228	1737963	45.662	nq	98
82) Chrysene	14.19	228	1692265	46.285	nq	99
83) Bis(2-ethylhexyl)phthalate	14.11	149	1248358	46.308	nq	99
84) Di-n-octyl phthalate	14.73	149	1896013	45.248	nq	99
85) Indeno(1,2,3-cd)pyrene	17.27	276	1402138	45.249	nq #	93
87) Benzo(b)fluoranthene	15.23	252	1441274	61.730	nq	98
88) Benzo(k)fluoranthene	15.27	252	1299779m	56.822	nq	
89) Benzo(a)pyrene	15.62	252	1241841	58.146	nq	98
90) Dibenzo(a,h)anthracene	17.28	278	1175057	63.648	nq #	96
91) Benzo(g,h,i)perylene	17.75	276	1184063	65.021	nq	93

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

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