

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120518\
 Data File : BF111105.D
 Acq On : 5 Dec 2018 13:27
 Operator : JU/SJ
 Sample : SSTDICC2.5
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC2.5

Manual Integrations
 APPROVED

mohammad
 12/6/2018 12:10:39 PM

Quant Time: Dec 05 17:25:01 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 05 15:25:21 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	476391	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	1985034	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	947588	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	1653218	20.00	ng	0.00
76) Chrysene-d12	13.95	240	1382272	20.00	ng	0.00
87) Perylene-d12	15.38	264	1255322	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.40	112	178409	6.21	ng	-0.01
7) Phenol-d6	6.42	99	228145	6.09	ng	-0.02
23) Nitrobenzene-d5	7.35	82	159035	4.59	ng	-0.02
42) 2,4,6-Tribromophenol	10.62	330	40331	4.29	ng	-0.01
45) 2-Fluorobiphenyl	0.00	172	0d	0.00	ng	
79) Terphenyl-d14	12.90	244	385520	5.43	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.52	88	46261	3.361	ng	# 80
3) Pyridine	3.28	79	96420	3.199	ng	# 69
4) n-Nitrosodimethylamine	3.18	42	47126	3.417	ng	# 81
6) Aniline	6.46	93	144969	3.132	ng	# 50
8) 2-Chlorophenol	6.58	128	102326	3.006	ng	97
10) Phenol	6.43	94	121305m	2.877	ng	
11) bis(2-Chloroethyl)ether	6.53	93	100324	3.159	ng	92
12) 1,3-Dichlorobenzene	6.75	146	111112	2.971	ng	97
13) 1,4-Dichlorobenzene	6.82	146	107846	2.804	ng	98
14) 1,2-Dichlorobenzene	6.97	146	104560	2.888	ng	96
15) Benzyl Alcohol	6.93	79	85507	2.972	ng	92
16) 2,2'-oxybis(1-Chloropropan	7.08	45	191797	3.794	ng	# 63
17) 2-Methylphenol	7.05	107	78267	2.891	ng	# 79
18) Hexachloroethane	7.32	117	38224	2.702	ng	93
19) n-Nitroso-di-n-propylamine	7.20	70	73029	2.995	ng	# 80
20) 3+4-Methylphenols	7.20	107	105601	2.935	ng	# 72
22) Acetophenone	7.20	105	146413	3.147	ng	# 95
24) Nitrobenzene	7.37	77	89417	2.479	ng	99
25) Isophorone	7.62	82	167731	2.948	ng	# 97
26) 2-Nitrophenol	7.69	139	27109	1.547	ng	97
27) 2,4-Dimethylphenol	7.73	122	80444	3.040	ng	96
28) bis(2-Chloroethoxy)methane	7.83	93	118048	3.097	ng	94
29) 2,4-Dichlorophenol	7.93	162	74043	2.674	ng	97
30) 1,2,4-Trichlorobenzene	8.03	180	80090	2.581	ng	94
33) 4-Chloroaniline	8.15	127	117472	2.966	ng	99
34) Hexachlorobutadiene	8.23	225	41951	2.373	ng	98
35) Caprolactam	8.46	113	27657	2.925	ng	# 79
36) 4-Chloro-3-methylphenol	8.62	107	83374	2.710	ng	99
37) 2-Methylnaphthalene	8.80	142	181137	2.947	ng	94
38) 1-Methylnaphthalene	8.89	142	169197	2.830	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.96	216	74772	2.508	ng	95
41) Hexachlorocyclopentadiene	8.96	237	32532	10.846	ng	95
43) 2,4,6-Trichlorophenol	9.07	196	49100	2.758	ng	98
44) 2,4,5-Trichlorophenol	9.10	196	46825	2.498	ng	# 91

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
47) 2-Chloronaphthalene	9.28	162	177413	2.801	nq	100
50) Dimethylphthalate	9.55	163	194117	2.793	nq	99
51) 2,6-Dinitrotoluene	9.61	165	25457	1.645	nq	97
52) Acenaphthene	9.87	154	168255	2.914	nq	98
53) 3-Nitroaniline	9.77	138	35145	1.949	nq #	95
56) 4-Nitrophenol	9.92	139	25592	2.846	nq #	82
59) 2,3,4,6-Tetrachlorophenol	10.15	232	35004	2.275	nq #	83
60) Diethylphthalate	10.25	149	190790	2.662	nq	99
62) 4-Nitroaniline	10.37	138	32784	1.953	nq	93
63) Azobenzene	10.53	77	209867	3.004	nq	95
66) n-Nitrosodiphenylamine	10.49	169	169650	3.083	nq	99
67) 4-Bromophenyl-phenylether	10.86	248	49359	2.741	nq #	84
68) Hexachlorobenzene	10.93	284	50788	2.648	nq #	80
69) Atrazine	11.01	200	51212	3.616	nq	96
70) Pentachlorophenol	11.12	266	26600	3.486	nq	97
77) Benzidine	12.64	184	160230	4.532	nq	97
78) Pyrene	12.75	202	277946	2.753	nq	96
80) Butylbenzylphthalate	13.38	149	110603	2.427	nq	99
81) Benzo(a)anthracene	13.93	228	218544	2.678	nq	99
82) 3,3'-Dichlorobenzidine	13.90	252	80147	2.864	nq #	98
83) Chrysene	13.97	228	216237	2.692	nq	99
84) Bis(2-ethylhexyl)phthalate	13.95	149	151678	2.497	nq #	96
85) Di-n-octyl phthalate	14.55	149	244580	2.551	nq #	92
86) Indeno(1,2,3-cd)pyrene	16.77	276	158955m	2.757	nq	
88) Benzo(b)fluoranthene	14.96	252	183790	2.492	nq #	97
89) Benzo(k)fluoranthene	14.99	252	194719	2.565	nq #	95
90) Benzo(a)pyrene	15.32	252	171212	2.494	nq	98
91) Dibenzo(a,h)anthracene	16.79	278	133378	2.355	nq #	90
92) Benzo(a,h,i)perylene	17.19	276	132752	2.431	nq #	84

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

