

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN073018\
 Data File : BN002188.D
 Acq On : 30 Jul 2018 16:19
 Operator : JU/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 ICVBN073018

Quant Time: Jul 30 16:49:10 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN073018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 30 16:17:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	3069	0.40	ng	0.00
7) Naphthalene-d8	10.47	136	12600	0.40	ng	0.00
13) Acenaphthene-d10	14.32	164	7109	0.40	ng	0.00
19) Phenanthrene-d10	17.07	188	14640	0.40	ng	0.00
27) Chrysene-d12	21.27	240	11019	0.40	ng	0.00
34) Perylene-d12	23.50	264	9911	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.32	112	2788	0.39	ng	0.00
5) Phenol-d6	6.88	99	3389	0.39	ng	0.00
8) Nitrobenzene-d5	8.84	82	3778	0.40	ng	0.00
11) 2-Methylnaphthalene-d10	12.06	152	7504	0.40	ng	0.00
14) 2,4,6-Tribromophenol	15.83	330	1051	0.35	ng	0.00
15) 2-Fluorobiphenyl	12.94	172	11088	0.41	ng	0.00
25) Fluoranthene-d10	19.11	212	13169	0.36	ng	0.00
29) Terphenyl-d14	19.71	244	9974	0.42	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.29	88	1340	0.388	ng	97
3) n-Nitrosodimethylamine	3.62	42	817	0.374	ng	# 88
6) bis(2-Chloroethyl)ether	7.13	93	3275	0.391	ng	89
9) Naphthalene	10.52	128	11657	0.410	ng	96
10) Hexachlorobutadiene	10.80	225	2521	0.410	ng	# 54
12) 2-Methylnaphthalene	12.13	142	7857	0.409	ng	98
16) Acenaphthylene	14.04	152	11920	0.392	ng	100
17) Acenaphthene	14.38	154	7857	0.402	ng	96
18) Fluorene	15.37	166	9638	0.401	ng	# 80
20) 4-Bromophenyl-phenylether	16.27	248	3406	0.411	ng	# 74
21) Hexachlorobenzene	16.38	284	3883	0.424	ng	96
22) Pentachlorophenol	16.74	266	554	0.335	ng	89
23) Phenanthrene	17.11	178	14744	0.405	ng	99
24) Anthracene	17.21	178	11553	0.383	ng	96
26) Fluoranthene	19.14	202	14406	0.363	ng	# 97
28) Pyrene	19.50	202	14379	0.441	ng	100
30) Benzo(a)anthracene	21.26	228	10874	0.392	ng	95
31) Chrysene	21.30	228	12792	0.412	ng	99
32) Bis(2-ethylhexyl)phthalate	21.18	149	4496	0.355	ng	# 94
33) Indeno(1,2,3-cd)pyrene	25.72	276	13093	0.448	ng	# 94
35) Benzo(b)fluoranthene	22.83	252	11473	0.394	ng	# 91
36) Benzo(k)fluoranthene	22.87	252	12120	0.401	ng	# 93
37) Benzo(a)pyrene	23.40	252	10716	0.406	ng	# 92
38) Dibenzo(a,h)anthracene	25.74	278	10834	0.422	ng	96
39) Benzo(g,h,i)perylene	26.40	276	11203	0.427	ng	# 90

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

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