

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN120518\
 Data File : BN003716.D
 Acq On : 05 Dec 2018 18:24
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.486

Quant Time: Dec 05 18:56:11 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 04 01:17:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	5086	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	19816	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	9954	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	22130	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	19960	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	17238	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	11761	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	24156	0.37	ng/ul	0.00

Target Compounds					Qvalue
3) Naphthalene	10.68	128	23068	0.414	ng/ul 100
5) 2-Methylnaphthalene	12.29	142	15595	0.401	ng/ul 100
7) Acenaphthylene	14.19	152	17812	0.397	ng/ul 100
8) Acenaphthene	14.53	153	16150	0.411	ng/ul 100
9) Fluorene	15.52	166	17998	0.393	ng/ul 98
11) Pentachlorophenol	16.86	266	1721	0.432	ng/ul 96
12) Phenanthrene	17.26	178	29849	0.411	ng/ul 100
13) Anthracene	17.35	178	24002	0.402	ng/ul 99
15) Fluoranthene	19.27	202	33329	0.389	ng/ul 99
17) Pyrene	19.63	202	32745	0.457	ng/ul 97
18) Benzo(a)anthracene	21.38	228	26039	0.413	ng/ul 99
19) Chrysene	21.43	228	29066	0.403	ng/ul 99
21) Benzo(b)fluoranthene	23.01	252	29757	0.417	ng/ul 99
22) Benzo(k)fluoranthene	23.05	252	28463	0.410	ng/ul 99
23) Benzo(a)pyrene	23.61	252	25765	0.417	ng/ul 99
24) Indeno(1,2,3-cd)pyrene	26.06	276	28180	0.398	ng/ul 98
25) Dibenzo(a,h)anthracene	26.07	278	22231	0.390	ng/ul 99
26) Benzo(g,h,i)perylene	26.78	276	24000	0.390	ng/ul 99

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

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