

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120618\
 Data File : BN003730.D
 Acq On : 06 Dec 2018 14:32
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.489

Quant Time: Dec 06 15:14:12 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 : Thu Dec 06 02:51:43 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	7212	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	29029	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	15158	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	33873	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	32378	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	29382	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	17680	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	41142	0.42	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.68	128	33887	0.416	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	23439	0.412	ng/ul	100
7) Acenaphthylene	14.19	152	29365	0.430	ng/ul	100
8) Acenaphthene	14.53	153	24825	0.415	ng/ul	100
9) Fluorene	15.52	166	27873	0.400	ng/ul	99
11) Pentachlorophenol	16.86	266	3013	0.494	ng/ul	99
12) Phenanthrene	17.25	178	46244	0.416	ng/ul	99
13) Anthracene	17.35	178	39447	0.432	ng/ul	100
15) Fluoranthene	19.27	202	55978	0.426	ng/ul	99
17) Pyrene	19.63	202	53888	0.464	ng/ul	98
18) Benzo(a)anthracene	21.38	228	46808	0.458	ng/ul	99
19) Chrysene	21.43	228	47627	0.407	ng/ul	100
21) Benzo(b)fluoranthene	23.01	252	48223	0.397	ng/ul	99
22) Benzo(k)fluoranthene	23.06	252	47398	0.400	ng/ul	99
23) Benzo(a)pyrene	23.61	252	43572	0.414	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.06	276	46942	0.389	ng/ul	98
25) Dibenzo(a,h)anthracene	26.07	278	36823	0.379	ng/ul	99
26) Benzo(g,h,i)perylene	26.78	276	39768	0.379	ng/ul	99

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

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