

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120518\
 Data File : BN003704.D
 Acq On : 05 Dec 2018 10:41
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.485

Quant Time: Dec 05 12:03:44 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	5368	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	21363	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	11328	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	25857	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	24565	0.40	ng/ul	0.00
20) Perylene-d12	23.70	264	22290	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	12971	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	31094	0.41	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.68	128	24764	0.413	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	17173	0.410	ng/ul	99
7) Acenaphthylene	14.20	152	21445	0.420	ng/ul	100
8) Acenaphthene	14.53	153	18351	0.411	ng/ul	100
9) Fluorene	15.52	166	20730	0.398	ng/ul	99
11) Pentachlorophenol	16.86	266	1930	0.414	ng/ul	98
12) Phenanthrene	17.26	178	35063	0.413	ng/ul	99
13) Anthracene	17.35	178	29598	0.424	ng/ul	100
15) Fluoranthene	19.27	202	42317	0.422	ng/ul	97
17) Pyrene	19.63	202	40647	0.461	ng/ul	99
18) Benzo(a)anthracene	21.38	228	34247	0.441	ng/ul	99
19) Chrysene	21.43	228	35938	0.405	ng/ul	100
21) Benzo(b)fluoranthene	23.00	252	36524	0.396	ng/ul	99
22) Benzo(k)fluoranthene	23.05	252	35990	0.401	ng/ul	99
23) Benzo(a)pyrene	23.60	252	32374	0.405	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.05	276	35283	0.385	ng/ul	97
25) Dibenzo(a,h)anthracene	26.07	278	27634	0.375	ng/ul	99
26) Benzo(g,h,i)perylene	26.77	276	30189	0.379	ng/ul	98

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

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