

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120418\
 Data File : BN003692.D
 Acq On : 04 Dec 2018 09:44
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.483

Quant Time: Dec 04 11:11:33 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	7009	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	27577	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	14697	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	33771	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	32811	0.40	ng/ul	0.00
20) Perylene-d12	23.70	264	29815	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	16401	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	40263	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	29936	0.386	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	20477	0.379	ng/ul	100
7) Acenaphthylene	14.19	152	28163	0.426	ng/ul	100
8) Acenaphthene	14.54	153	22980	0.396	ng/ul	99
9) Fluorene	15.52	166	25243	0.373	ng/ul	99
11) Pentachlorophenol	16.86	266	3032	0.498	ng/ul	99
12) Phenanthrene	17.26	178	43055	0.388	ng/ul	100
13) Anthracene	17.35	178	35323	0.388	ng/ul	99
15) Fluoranthene	19.27	202	52054	0.398	ng/ul	99
17) Pyrene	19.63	202	51659	0.439	ng/ul	98
18) Benzo(a)anthracene	21.38	228	42602	0.411	ng/ul	99
19) Chrysene	21.43	228	45436	0.383	ng/ul	100
21) Benzo(b)fluoranthene	23.00	252	46417	0.376	ng/ul	100
22) Benzo(k)fluoranthene	23.05	252	44930	0.374	ng/ul	99
23) Benzo(a)pyrene	23.60	252	40132	0.376	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.05	276	43119	0.352	ng/ul	98
25) Dibenzo(a,h)anthracene	26.07	278	34030	0.345	ng/ul	99
26) Benzo(g,h,i)perylene	26.78	276	36489	0.343	ng/ul	99

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

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