

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120618\
 Data File : BN003743.D
 Acq On : 06 Dec 2018 22:20
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.490

Quant Time: Dec 07 03:20:52 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 : Fri Dec 07 03:02:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	4862	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	19350	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	9811	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	21427	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	18368	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	15568	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	11731	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	23613	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	22493	0.414	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	15555	0.410	ng/ul	99
7) Acenaphthylene	14.19	152	18587	0.421	ng/ul	100
8) Acenaphthene	14.53	153	16134	0.417	ng/ul	100
9) Fluorene	15.52	166	17855	0.396	ng/ul	99
11) Pentachlorophenol	16.86	266	1550	0.402	ng/ul	100
12) Phenanthrene	17.25	178	29272	0.416	ng/ul	99
13) Anthracene	17.35	178	24354	0.421	ng/ul	100
15) Fluoranthene	19.27	202	32316	0.389	ng/ul	96
17) Pyrene	19.63	202	31427	0.477	ng/ul	99
18) Benzo(a)anthracene	21.38	228	25066	0.432	ng/ul	98
19) Chrysene	21.43	228	27648	0.416	ng/ul	100
21) Benzo(b)fluoranthene	23.01	252	27315	0.424	ng/ul	98
22) Benzo(k)fluoranthene	23.05	252	26370	0.420	ng/ul	99
23) Benzo(a)pyrene	23.61	252	23635	0.424	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.06	276	25831	0.404	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.07	278	20590	0.400	ng/ul	99
26) Benzo(g,h,i)perylene	26.78	276	21841	0.393	ng/ul	98

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

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