

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121118\
 Data File : BN003806.D
 Acq On : 11 Dec 2018 09:30
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.497

Quant Time: Dec 11 14:55:50 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 11 04:22:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	4872	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	19817	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	9691	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	19153	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	13987	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	14055	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	11130	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	18297	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	21265	0.382	ng/ul	99
5) 2-Methylnaphthalene	12.29	142	14587	0.375	ng/ul	99
7) Acenaphthylene	14.19	152	19340	0.443	ng/ul	100
8) Acenaphthene	14.53	153	14661	0.383	ng/ul	98
9) Fluorene	15.52	166	15857	0.356	ng/ul	99
11) Pentachlorophenol	16.86	266	1356	0.393	ng/ul	99
12) Phenanthrene	17.26	178	23638	0.376	ng/ul	100
13) Anthracene	17.35	178	21866	0.423	ng/ul	100
15) Fluoranthene	19.27	202	24334	0.328	ng/ul	97
17) Pyrene	19.64	202	24108	0.481	ng/ul	99
18) Benzo(a)anthracene	21.39	228	19881	0.450	ng/ul	98
19) Chrysene	21.43	228	19233	0.380	ng/ul	98
21) Benzo(b)fluoranthene	23.01	252	19906	0.342	ng/ul	91
22) Benzo(k)fluoranthene	23.06	252	19846	0.350	ng/ul#	92
23) Benzo(a)pyrene	23.61	252	19402	0.385	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	26.07	276	23903	0.414	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.08	278	19009	0.409	ng/ul#	93
26) Benzo(g,h,i)perylene	26.79	276	20195	0.402	ng/ul	94

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

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