

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121420\
 Data File : BN013107.D
 Acq On : 14 Dec 2020 09:12
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.442

Quant Time: Dec 14 10:25:13 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 14 10:24:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.36	152	1131	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	4144	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	2318	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.79	188	4997	0.40	ng/ul	0.00
16) Chrysene-d12	21.02	240	4100	0.40	ng/ul	0.00
20) Perylene-d12	23.12	264	4474	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	2589	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.84	212	5791	0.43	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.17	128	5061	0.422	ng/ul	93
5) 2-Methylnaphthalene	11.80	142	3356	0.426	ng/ul	100
7) Acenaphthylene	13.75	152	5154	0.470	ng/ul	96
8) Acenaphthene	14.09	153	3908	0.434	ng/ul	99
9) Fluorene	15.09	166	4429	0.429	ng/ul	99
11) Pentachlorophenol	16.46	266	707	0.488	ng/ul	94
12) Phenanthrene	16.84	178	7231	0.432	ng/ul	99
13) Anthracene	16.92	178	6847	0.480	ng/ul	97
15) Fluoranthene	18.87	202	8451	0.424	ng/ul	100
17) Pyrene	19.24	202	8922	0.461	ng/ul	97
18) Benzo(a)anthracene	21.00	228	7500	0.442	ng/ul	99
19) Chrysene	21.06	228	7594	0.405	ng/ul	98
21) Benzo(b)fluoranthene	22.50	252	8219	0.397	ng/ul	94
22) Benzo(k)fluoranthene	22.54	252	8713	0.381	ng/ul#	96
23) Benzo(a)pyrene	23.03	252	7955	0.414	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.17	276	9698	0.384	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.18	278	7726	0.387	ng/ul#	95
26) Benzo(g,h,i)perylene	25.79	276	8308	0.364	ng/ul	96

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

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