

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN010220\
 Data File : BN009287.D
 Acq On : 02 Jan 2020 10:10
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.409

Quant Time: Jan 02 12:44:47 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN121619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 02 12:41:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	2208	0.40	ng/ul	0.00
2) Naphthalene-d8	10.26	136	7764	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.13	164	3889	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.88	188	7177	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	5475	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	5526	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.85	152	4910	0.36	ng/ul	0.00
14) Fluoranthene-d10	18.92	212	8194	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.30	128	9526	0.405	ng/ul	99
5) 2-Methylnaphthalene	11.93	142	6337	0.356	ng/ul	100
7) Acenaphthylene	13.85	152	7926	0.388	ng/ul	98
8) Acenaphthene	14.19	153	6195	0.386	ng/ul	99
9) Fluorene	15.19	166	6616	0.349	ng/ul	100
11) Pentachlorophenol	16.55	266	782	0.409	ng/ul	98
12) Phenanthrene	16.92	178	10086	0.404	ng/ul	100
13) Anthracene	17.01	178	8978	0.401	ng/ul	99
15) Fluoranthene	18.95	202	10906	0.382	ng/ul	99
17) Pyrene	19.31	202	11034	0.395	ng/ul	99
18) Benzo(a)anthracene	21.07	228	9983	0.411	ng/ul	99
19) Chrysene	21.12	228	10447	0.428	ng/ul	99
21) Benzo(b)fluoranthene	22.58	252	9958	0.386	ng/ul	98
22) Benzo(k)fluoranthene	22.62	252	10476	0.403	ng/ul	97
23) Benzo(a)pyrene	23.12	252	9336	0.413	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.30	276	10908	0.443	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.31	278	8736	0.446	ng/ul	98
26) Benzo(g,h,i)perylene	25.93	276	9513	0.459	ng/ul	98

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

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