

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082819\
 Data File : BN007442.D
 Acq On : 27 Aug 2019 21:17
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.433

Quant Time: Aug 27 21:48:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN082819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 27 20:30:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	3968	0.40	ng/ul	0.00
2) Naphthalene-d8	10.16	136	16152	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	8178	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	18614	0.40	ng/ul	0.00
16) Chrysene-d12	21.02	240	16143	0.40	ng/ul	0.00
20) Perylene-d12	23.12	264	17037	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	9754	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.84	212	23217	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.21	128	17774	0.390	ng/ul	99
5) 2-Methylnaphthalene	11.84	142	11845	0.387	ng/ul	99
7) Acenaphthylene	13.76	152	17592	0.387	ng/ul	99
8) Acenaphthene	14.11	153	12778	0.391	ng/ul	99
9) Fluorene	15.11	166	13651	0.386	ng/ul	99
11) Pentachlorophenol	16.47	266	1411	0.336	ng/ul	98
12) Phenanthrene	16.84	178	22597	0.389	ng/ul	100
13) Anthracene	16.94	178	21198	0.387	ng/ul	100
15) Fluoranthene	18.87	202	29500	0.379	ng/ul	99
17) Pyrene	19.24	202	30706	0.389	ng/ul	99
18) Benzo(a)anthracene	21.01	228	27802	0.385	ng/ul	99
19) Chrysene	21.06	228	30331	0.391	ng/ul	100
21) Benzo(b)fluoranthene	22.51	252	28928	0.389	ng/ul	99
22) Benzo(k)fluoranthene	22.54	252	30628	0.394	ng/ul	97
23) Benzo(a)pyrene	23.04	252	28364	0.382	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.19	276	32101	0.387	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.20	278	25404	0.387	ng/ul	99
26) Benzo(g,h,i)perylene	25.81	276	27455	0.387	ng/ul	100

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

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