

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071819\
 Data File : BN006906.D
 Acq On : 18 Jul 2019 23:39
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.436

Quant Time: Jul 19 01:15:11 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN062719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 06:39:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	2279	0.40	ng/ul	0.01
2) Naphthalene-d8	10.37	136	8823	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	4667	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	10114	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	8174	0.40	ng/ul	-0.01
20) Perylene-d12	23.45	264	9094	0.40	ng/ul	-0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	5052	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	10778	0.36	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.42	128	9956	0.393	ng/ul	98
5) 2-Methylnaphthalene	12.06	142	6049	0.353	ng/ul	98
7) Acenaphthylene	13.96	152	8729	0.362	ng/ul	99
8) Acenaphthene	14.30	153	7096	0.382	ng/ul	98
9) Fluorene	15.31	166	7376	0.363	ng/ul	97
11) Pentachlorophenol	16.72	266	443	0.294	ng/ul	96
12) Phenanthrene	17.05	178	12007	0.369	ng/ul	99
13) Anthracene	17.15	178	9440	0.336	ng/ul	99
15) Fluoranthene	19.08	202	14037	0.349	ng/ul	98
17) Pyrene	19.45	202	14608	0.391	ng/ul	99
18) Benzo(a)anthracene	21.21	228	11572	0.345	ng/ul	100
19) Chrysene	21.26	228	15090	0.403	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	12974	0.348	ng/ul	98
22) Benzo(k)fluoranthene	22.83	252	15203	0.390	ng/ul	98
23) Benzo(a)pyrene	23.36	252	13623	0.373	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.68	276	15643	0.423	ng/ul	94
25) Dibenzo(a,h)anthracene	25.68	278	12199	0.415	ng/ul	97
26) Benzo(g,h,i)perylene	26.36	276	13761	0.445	ng/ul	97

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

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