

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071819\
 Data File : BN006894.D
 Acq On : 18 Jul 2019 16:42
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.435

Quant Time: Jul 18 19:01:24 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	2150	0.40	ng/ul	0.01
2) Naphthalene-d8	10.37	136	8430	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	4422	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	9246	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	7519	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	8324	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	4735	0.37	ng/ul	0.01
14) Fluoranthene-d10	19.05	212	9844	0.36	ng/ul	0.00

Target Compounds						Ovalue
3) Naphthalene	10.42	128	9570	0.395	ng/ul	97
5) 2-Methylnaphthalene	12.06	142	5775	0.353	ng/ul	99
7) Acenaphthylene	12.97	152	8234	0.361	ng/ul	99
8) Acenaphthene	14.31	153	6696	0.381	ng/ul	99
9) Fluorene	15.31	166	6833	0.354	ng/ul	98
11) Pentachlorophenol	16.72	266	399	0.290	ng/ul	97
12) Phenanthrene	17.05	178	11180	0.376	ng/ul	99
13) Anthracene	17.16	178	8545	0.332	ng/ul	98
15) Fluoranthene	19.09	202	12735	0.346	ng/ul	99
17) Pyrene	19.45	202	13493	0.393	ng/ul	97
18) Benzo(a)anthracene	21.22	228	10592	0.343	ng/ul	99
19) Chrysene	21.27	228	13907	0.403	ng/ul	99
21) Benzo(b)fluoranthene	22.80	252	12169	0.357	ng/ul	97
22) Benzo(k)fluoranthene	22.84	252	14066	0.394	ng/ul	98
23) Benzo(a)pyrene	23.37	252	12380	0.370	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.69	276	14507	0.429	ng/ul	93
25) Dibenzo(a,h)anthracene	25.69	278	11298	0.420	ng/ul	97
26) Benzo(g,h,i)perylene	26.37	276	12757	0.451	ng/ul	97

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

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