

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN060819\
 Data File : BN006183.D
 Acq On : 08 Jun 2019 10:06
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 DFTPP87

Quant Time: Jun 08 11:21:27 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 07 01:31:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	5475	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	21024	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	9673	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	18766	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	14938	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	16835	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	11447	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	19517	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	22456	0.387	ng/ul	99
5) 2-Methylnaphthalene	12.10	142	14286	0.353	ng/ul	98
7) Acenaphthylene	14.01	152	20100	0.374	ng/ul	100
8) Acenaphthene	14.36	153	14555	0.390	ng/ul	99
9) Fluorene	15.35	166	15098	0.346	ng/ul	97
11) Pentachlorophenol	16.71	266	1561	0.327	ng/ul	96
12) Phenanthrene	17.09	178	22710	0.377	ng/ul	100
13) Anthracene	17.18	178	20122	0.350	ng/ul	100
15) Fluoranthene	19.12	202	26138	0.393	ng/ul	98
17) Pyrene	19.49	202	27244	0.346	ng/ul	98
18) Benzo(a)anthracene	21.25	228	23201	0.351	ng/ul	98
19) Chrysene	21.31	228	26832	0.433	ng/ul	99
21) Benzo(b)fluoranthene	22.84	252	24962	0.356	ng/ul	97
22) Benzo(k)fluoranthene	22.88	252	27245	0.398	ng/ul	95
23) Benzo(a)pyrene	23.41	252	25653	0.392	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.74	276	29477	0.417	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.75	278	23462	0.415	ng/ul	93
26) Benzo(g,h,i)perylene	26.42	276	24900	0.420	ng/ul	95

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

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