

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM073119\
 Data File : BM021759.D
 Acq On : 01 Aug 2019 00:10
 Operator : HP/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.451

Quant Time: Aug 01 01:25:34 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM073119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 31 15:40:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	2114	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	4794	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.32	164	1714	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.08	188	4218	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	4479	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	5486	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	2150	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.11	212	4556	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.51	128	4670	0.386	ng/ul	99
5) 2-Methylnaphthalene	12.14	142	2409	0.377	ng/ul	93
7) Acenaphthylene	14.04	152	3644	0.389	ng/ul	98
8) Acenaphthene	14.39	153	2640	0.382	ng/ul	97
9) Fluorene	15.38	166	2953	0.381	ng/ul	99
11) Pentachlorophenol	16.75	266	559	0.388	ng/ul	98
12) Phenanthrene	17.12	178	4562	0.391	ng/ul	98
13) Anthracene	17.22	178	3988	0.397	ng/ul	99
15) Fluoranthene	19.14	202	5374	0.369	ng/ul	100
17) Pyrene	19.51	202	5404	0.426	ng/ul	99
18) Benzo(a)anthracene	21.26	228	5542	0.397	ng/ul	99
19) Chrysene	21.31	228	6738	0.383	ng/ul	99
21) Benzo(b)fluoranthene	22.83	252	7934	0.392	ng/ul	96
22) Benzo(k)fluoranthene	22.88	252	7986	0.389	ng/ul	97
23) Benzo(a)pyrene	23.41	252	6994	0.384	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.75	276	8613	0.354	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.76	278	6786	0.355	ng/ul	97
26) Benzo(g,h,i)perylene	26.44	276	8128	0.348	ng/ul	96

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

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