

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM073119\
 Data File : BM021754.D
 Acq On : 31 Jul 2019 21:08
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.499

Quant Time: Jul 31 21:40:11 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 20:59:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	1957	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	4489	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.32	164	1561	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.08	188	3950	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	4695	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	6398	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	1990	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.11	212	4235	0.37	ng/ul	0.00

Target Compounds					Qvalue
3) Naphthalene	10.51	128	4326	0.382	ng/ul 99
5) 2-Methylnaphthalene	12.14	142	2187	0.365	ng/ul 98
7) Acenaphthylene	14.04	152	3364	0.394	ng/ul 99
8) Acenaphthene	14.39	153	2453	0.390	ng/ul 99
9) Fluorene	15.39	166	2680	0.380	ng/ul 100
11) Pentachlorophenol	16.75	266	504	0.374	ng/ul 99
12) Phenanthrene	17.12	178	4209	0.385	ng/ul 99
13) Anthracene	17.22	178	3700	0.393	ng/ul 100
15) Fluoranthene	19.14	202	5009	0.367	ng/ul 100
17) Pyrene	19.51	202	5158	0.388	ng/ul 100
18) Benzo(a)anthracene	21.26	228	5621	0.384	ng/ul 100
19) Chrysene	21.31	228	7198	0.390	ng/ul 98
21) Benzo(b)fluoranthene	22.84	252	9007	0.381	ng/ul 99
22) Benzo(k)fluoranthene	22.88	252	8952	0.374	ng/ul 98
23) Benzo(a)pyrene	23.42	252	7893	0.372	ng/ul 96
24) Indeno(1,2,3-cd)pyrene	25.75	276	9502	0.334	ng/ul# 97
25) Dibenzo(a,h)anthracene	25.77	278	7448	0.335	ng/ul 95
26) Benzo(g,h,i)perylene	26.45	276	8960	0.329	ng/ul 97

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

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