

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040420\
 Data File : BM025930.D
 Acq On : 04 Apr 2020 02:00
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.438

Quant Time: Apr 04 02:45:04 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 03 22:53:21 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	2341	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	9940	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	6200	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	15976	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	16565	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	17267	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	6799	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	20303	0.41	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.37	128	10598	0.363	ng/ul	100
5) 2-Methylnaphthalene	12.00	142	7801	0.381	ng/ul	99
7) Acenaphthylene	13.92	152	11367	0.429	ng/ul	99
8) Acenaphthene	14.26	153	8771	0.399	ng/ul	99
9) Fluorene	15.26	166	10858	0.398	ng/ul	100
11) Pentachlorophenol	16.61	266	1458	0.372	ng/ul	99
12) Phenanthrene	16.99	178	18805	0.374	ng/ul	100
13) Anthracene	17.08	178	16851	0.379	ng/ul	100
15) Fluoranthene	19.01	202	23315	0.375	ng/ul	99
17) Pyrene	19.37	202	23619	0.398	ng/ul	96
18) Benzo(a)anthracene	21.12	228	23013	0.389	ng/ul	100
19) Chrysene	21.17	228	24586	0.380	ng/ul	100
21) Benzo(b)fluoranthene	22.65	252	25272	0.398	ng/ul	99
22) Benzo(k)fluoranthene	22.69	252	25890	0.380	ng/ul	99
23) Benzo(a)pyrene	23.19	252	23511	0.414	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.40	276	27234	0.358	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.42	278	22676	0.357	ng/ul	99
26) Benzo(g,h,i)perylene	26.04	276	22384	0.348	ng/ul	98

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

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