

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM050120\
 Data File : BM025964.D
 Acq On : 01 May 2020 11:11
 Operator : MA/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.483

Quant Time: May 01 12:26:05 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 01 11:47:12 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	1367	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	5112	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	3167	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	7291	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	5256	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	4338	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	2885	0.34	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	6607	0.32	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.37	128	5399	0.34	ng/ul	100
5) 2-Methylnaphthalene	11.99	142	3594	0.34	ng/ul	98
7) Acenaphthylene	13.91	152	4683	0.32	ng/ul	99
8) Acenaphthene	14.26	153	4065	0.33	ng/ul	99
9) Fluorene	15.25	166	4737	0.32	ng/ul	100
11) Pentachlorophenol	16.61	266	580	0.31	ng/ul	92
12) Phenanthrene	16.99	178	8054	0.32	ng/ul	99
13) Anthracene	17.08	178	6686	0.32	ng/ul	99
15) Fluoranthene	19.00	202	9045	0.31	ng/ul	98
17) Pyrene	19.37	202	9013	0.36	ng/ul	99
18) Benzo(a)anthracene	21.13	228	6416	0.32	ng/ul	99
19) Chrysene	21.17	228	7915	0.32	ng/ul	100
21) Benzo(b)fluoranthene	22.65	252	6295	0.32	ng/ul	99
22) Benzo(k)fluoranthene	22.69	252	6838	0.31	ng/ul	99
23) Benzo(a)pyrene	23.20	252	5146	0.31	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.41	276	5635	0.31	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.42	278	4377	0.31	ng/ul	97
26) Benzo(g,h,i)perylene	26.05	276	5302	0.32	ng/ul	98

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

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