

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM050420\
 Data File : BM025970.D
 Acq On : 04 May 2020 11:35
 Operator : MA/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.485

Quant Time: May 04 12:06:24 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 12:24:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	1633	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	5531	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	3166	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	7300	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	6720	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	6421	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	3530	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	8483	0.41	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.36	128	7001	0.40	ng/ul	98
5) 2-Methylnaphthalene	11.99	142	4481	0.39	ng/ul	99
7) Acenaphthylene	13.91	152	5613	0.38	ng/ul	99
8) Acenaphthene	14.25	153	4911	0.39	ng/ul	98
9) Fluorene	15.25	166	5798	0.40	ng/ul	99
11) Pentachlorophenol	16.61	266	735	0.40	ng/ul	99
12) Phenanthrene	16.98	178	9791	0.39	ng/ul	100
13) Anthracene	17.07	178	8326	0.40	ng/ul	99
15) Fluoranthene	19.00	202	11623	0.40	ng/ul	98
17) Pyrene	19.37	202	11704	0.36	ng/ul	99
18) Benzo(a)anthracene	21.12	228	10345	0.41	ng/ul	100
19) Chrysene	21.17	228	12016	0.38	ng/ul	99
21) Benzo(b)fluoranthene	22.65	252	10868	0.37	ng/ul	97
22) Benzo(k)fluoranthene	22.69	252	12109	0.38	ng/ul	97
23) Benzo(a)pyrene	23.19	252	9629	0.39	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.40	276	12062	0.45	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.41	278	9921	0.48	ng/ul	92
26) Benzo(g,h,i)perylene	26.04	276	10355	0.42	ng/ul	96

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

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