

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025206.D
 Acq On : 03 Mar 2020 10:43
 Operator : CG/JU
 Sample : SSTD0.178
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.178

Quant Time: Mar 03 12:39:22 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 03 12:35:16 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	2642	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	9672	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	6542	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	16415	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	16460	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	16342	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	1761	0.10	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	5253	0.09	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.50	128	2924	0.099	ng/ul#	87
5) 2-Methylnaphthalene	12.12	142	2090	0.098	ng/ul	98
7) Acenaphthylene	14.02	152	3094	0.082	ng/ul	99
8) Acenaphthene	14.37	153	2389	0.094	ng/ul	97
9) Fluorene	15.36	166	2900	0.091	ng/ul	97
12) Phenanthrene	17.09	178	5221	0.090	ng/ul	99
13) Anthracene	17.18	178	4424	0.079	ng/ul	98
15) Fluoranthene	19.11	202	6300	0.075	ng/ul	96
17) Pyrene	19.47	202	6407	0.086	ng/ul	96
18) Benzo(a)anthracene	21.22	228	5892	0.083	ng/ul	98
19) Chrysene	21.27	228	6720	0.095	ng/ul	98
21) Benzo(b)fluoranthene	22.77	252	6294	0.098	ng/ul	94
22) Benzo(k)fluoranthene	22.82	252	6640	0.105	ng/ul	93
23) Benzo(a)pyrene	23.33	252	5216	0.089	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.61	276	6456	0.088	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.62	278	5155	0.084	ng/ul#	91
26) Benzo(a,h,i)perylene	26.27	276	5482	0.089	ng/ul#	94

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

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