

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025287.D
 Acq On : 05 Mar 2020 14:38
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 83 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.435

Quant Time: Mar 05 15:24:46 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 : Wed Mar 04 08:02:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	3467	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	14054	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	10047	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	25139	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	23354	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	19521	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	9305	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	28314	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.49	128	15034	0.355	ng/ul#	83
5) 2-Methylnaphthalene	12.11	142	10961	0.356	ng/ul	99
7) Acenaphthylene	14.02	152	16397	0.349	ng/ul#	93
8) Acenaphthene	14.36	153	13089	0.359	ng/ul	99
9) Fluorene	15.35	166	16340	0.357	ng/ul	95
11) Pentachlorophenol	16.70	266	2229	0.339	ng/ul	98
12) Phenanthrene	17.08	178	28694	0.354	ng/ul	95
13) Anthracene	17.17	178	25664	0.357	ng/ul	96
15) Fluoranthene	19.10	202	35025	0.345	ng/ul	94
17) Pyrene	19.46	202	34782	0.393	ng/ul#	94
18) Benzo(a)anthracene	21.21	228	31368	0.365	ng/ul	98
19) Chrysene	21.26	228	33748	0.358	ng/ul	98
21) Benzo(b)fluoranthene	22.76	252	28271	0.362	ng/ul	93
22) Benzo(k)fluoranthene	22.81	252	31272	0.382	ng/ul	98
23) Benzo(a)pyrene	23.32	252	23750	0.362	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.60	276	28975	0.355	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.61	278	24069	0.355	ng/ul#	94
26) Benzo(g,h,i)perylene	26.26	276	23281	0.335	ng/ul#	92

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

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