

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM111020\
 Data File : BM027771.D
 Acq On : 10 Nov 2020 15:12
 Operator : JU/CG
 Sample : SSTD3.277
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.277

Quant Time: Nov 10 15:43:12 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM111020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 10 14:33:58 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.39	152	545	0.40	ng/ul	0.00
2) Naphthalene-d8	11.22	136	1970	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.00	164	1257	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	2734	0.40	ng/ul	0.00
16) Chrysene-d12	21.81	240	2613	0.40	ng/ul	0.00
20) Perylene-d12	24.25	264	2365	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.78	152	10950	3.02	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	25987	3.19	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.27	128	18002	2.859	ng/ul	95
5) 2-Methylnaphthalene	12.86	142	12861	2.962	ng/ul	100
7) Acenaphthylene	14.72	152	18878	3.019	ng/ul	96
8) Acenaphthene	15.06	153	14105	2.975	ng/ul	99
9) Fluorene	16.03	166	17202	3.071	ng/ul	99
11) Pentachlorophenol	17.35	266	3138	2.825	ng/ul	97
12) Phenanthrene	17.75	178	26782	3.020	ng/ul	96
13) Anthracene	17.84	178	27337	3.186	ng/ul	98
15) Fluoranthene	19.73	202	33027	3.269	ng/ul	97
17) Pyrene	20.08	202	33104	2.902	ng/ul	98
18) Benzo(a)anthracene	21.79	228	31113	3.007	ng/ul	98
19) Chrysene	21.85	228	31854	3.012	ng/ul	97
21) Benzo(b)fluoranthene	23.50	252	31340	3.013	ng/ul	93
22) Benzo(k)fluoranthene	23.55	252	31662	3.059	ng/ul#	94
23) Benzo(a)pyrene	24.14	252	28805	3.141	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	26.77	276	34172	3.147	ng/ul#	96
25) Dibenzo(a,h)anthracene	26.78	278	28696	3.175	ng/ul	95
26) Benzo(g,h,i)perylene	27.55	276	29068	3.104	ng/ul	95

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

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