

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092320\
 Data File : BM027498.D
 Acq On : 23 Sep 2020 11:57
 Operator : JU/CG
 Sample : SSTD0.879
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.879

Quant Time: Sep 23 12:28:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM092320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 23 11:54:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	1305	0.40	ng/ul	0.00
2) Naphthalene-d8	10.61	136	3935	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.45	164	2745	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.20	188	6212	0.40	ng/ul	0.00
16) Chrysene-d12	21.37	240	5061	0.40	ng/ul	0.00
20) Perylene-d12	23.65	264	4864	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.20	152	6067	0.91	ng/ul	0.00
14) Fluoranthene-d10	19.22	212	14276	0.81	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.66	128	9347	0.793	ng/ul	98
5) 2-Methylnaphthalene	12.27	142	6889	0.854	ng/ul	100
7) Acenaphthylene	14.17	152	11600	0.973	ng/ul	98
8) Acenaphthene	14.51	153	8144	0.786	ng/ul	98
9) Fluorene	15.50	166	10033	0.808	ng/ul	100
11) Pentachlorophenol	16.84	266	1360	0.664	ng/ul	97
12) Phenanthrene	17.24	178	15833	0.754	ng/ul	99
13) Anthracene	17.33	178	15583	0.817	ng/ul	99
15) Fluoranthene	19.25	202	19012	0.753	ng/ul	99
17) Pyrene	19.61	202	18693	0.864	ng/ul	98
18) Benzo(a)anthracene	21.36	228	16257	0.809	ng/ul	100
19) Chrysene	21.41	228	16757	0.770	ng/ul	100
21) Benzo(b)fluoranthene	22.97	252	16276	0.780	ng/ul	98
22) Benzo(k)fluoranthene	23.01	252	17131	0.790	ng/ul#	97
23) Benzo(a)pyrene	23.55	252	14737	0.780	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.95	276	19283	0.740	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.97	278	15945	0.752	ng/ul	98
26) Benzo(g,h,i)perylene	26.65	276	16290	0.724	ng/ul	98

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

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