

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092320\
 Data File : BM027499.D
 Acq On : 23 Sep 2020 12:33
 Operator : JU/CG
 Sample : SSTD1.680
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD1.680

Quant Time: Sep 23 13:28:00 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	1326	0.40	ng/ul	0.00
2) Naphthalene-d8	10.61	136	4124	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.45	164	3056	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.19	188	7321	0.40	ng/ul	0.00
16) Chrysene-d12	21.37	240	5552	0.40	ng/ul	0.00
20) Perylene-d12	23.65	264	4859	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.20	152	12689	1.82	ng/ul	0.00
14) Fluoranthene-d10	19.22	212	33203	1.59	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.66	128	19075	1.545	ng/ul	97
5) 2-Methylnaphthalene	12.27	142	14472	1.711	ng/ul	99
7) Acenaphthylene	14.17	152	25124	1.893	ng/ul	98
8) Acenaphthene	14.51	153	17793	1.543	ng/ul	97
9) Fluorene	15.50	166	22619	1.637	ng/ul	98
11) Pentachlorophenol	16.84	266	3354	1.389	ng/ul	100
12) Phenanthrene	17.23	178	36707	1.483	ng/ul	99
13) Anthracene	17.32	178	36649	1.630	ng/ul	99
15) Fluoranthene	19.25	202	44207	1.486	ng/ul	98
17) Pyrene	19.61	202	43070	1.815	ng/ul	98
18) Benzo(a)anthracene	21.36	228	35152	1.594	ng/ul	100
19) Chrysene	21.41	228	35901	1.504	ng/ul	100
21) Benzo(b)fluoranthene	22.97	252	33436	1.605	ng/ul	97
22) Benzo(k)fluoranthene	23.01	252	34085	1.573	ng/ul#	96
23) Benzo(a)pyrene	23.55	252	29560	1.566	ng/ul#	96
24) Indeno(1,2,3-cd)pyrene	25.95	276	37878	1.455	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.97	278	31273	1.476	ng/ul#	96
26) Benzo(g,h,i)perylene	26.65	276	31831	1.417	ng/ul	98

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

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