

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041520\
 Data File : BM025959.D
 Acq On : 15 Apr 2020 10:39
 Operator : MA/SJ
 Sample : SSTD0.877
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.877

Quant Time: Apr 15 11:19:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 15 11:17:24 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	946	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	3485	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	2152	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	4922	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	4326	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	3641	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	4903	0.79	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	11927	0.77	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.37	128	9178	0.896	ng/ul	97
5) 2-Methylnaphthalene	11.99	142	6205	0.865	ng/ul	98
7) Acenaphthylene	13.91	152	8447	0.918	ng/ul	99
8) Acenaphthene	14.26	153	7061	0.925	ng/ul	99
9) Fluorene	15.25	166	8477	0.895	ng/ul	99
11) Pentachlorophenol	16.61	266	1041	0.862	ng/ul	97
12) Phenanthrene	16.98	178	13993	0.902	ng/ul	98
13) Anthracene	17.07	178	12160	0.888	ng/ul	98
15) Fluoranthene	19.00	202	16598	0.866	ng/ul	99
17) Pyrene	19.37	202	16434	1.061	ng/ul	98
18) Benzo(a)anthracene	21.12	228	13852	0.896	ng/ul	99
19) Chrysene	21.17	228	16600	0.984	ng/ul	99
21) Benzo(b)fluoranthene	22.65	252	13909	1.038	ng/ul	96
22) Benzo(k)fluoranthene	22.69	252	15291	1.064	ng/ul	94
23) Benzo(a)pyrene	23.19	252	11660	0.974	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.40	276	13062	0.814	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.42	278	10240	0.765	ng/ul	92
26) Benzo(g,h,i)perylene	26.04	276	11863	0.874	ng/ul	95

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

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