

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081720\
 Data File : BM027113.D
 Acq On : 17 Aug 2020 13:05
 Operator : JU/CG
 Sample : SSTD1.679
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD1.679

Quant Time: Aug 17 13:40:54 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM081720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 17 12:32:35 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1227	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	3857	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	2457	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	5481	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	5464	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	5416	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	10828	1.71	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	25893	1.65	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.45	128	18089	1.490	ng/ul	98
5) 2-Methylnaphthalene	12.07	142	12775	1.590	ng/ul	96
7) Acenaphthylene	13.98	152	16595	1.449	ng/ul	97
8) Acenaphthene	14.33	153	14560	1.503	ng/ul	99
9) Fluorene	15.32	166	17944	1.581	ng/ul	98
11) Pentachlorophenol	16.67	266	2912	2.103	ng/ul	97
12) Phenanthrene	17.05	178	29560	1.581	ng/ul	98
13) Anthracene	17.14	178	27400	1.741	ng/ul	97
15) Fluoranthene	19.07	202	36491	1.678	ng/ul	98
17) Pyrene	19.43	202	36265	1.377	ng/ul	99
18) Benzo(a)anthracene	21.19	228	34771	1.682	ng/ul	98
19) Chrysene	21.24	228	37037	1.442	ng/ul	99
21) Benzo(b)fluoranthene	22.75	252	37841	1.530	ng/ul	92
22) Benzo(k)fluoranthene	22.79	252	39431	1.449	ng/ul#	92
23) Benzo(a)pyrene	23.30	252	34239	1.652	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.59	276	47865	2.096	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.60	278	39762	2.266	ng/ul	93
26) Benzo(g,h,i)perylene	26.25	276	41411	1.980	ng/ul	96

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

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