

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025210.D
 Acq On : 03 Mar 2020 13:09
 Operator : CG/JU
 Sample : SSTD1.676
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD1.676

Quant Time: Mar 03 13:47:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 03 12:35:16 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	2162	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	8273	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	6099	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	14485	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	15769	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	13646	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	26230	1.69	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	80996	1.63	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.50	128	40993	1.628	ng/ul#	82
5) 2-Methylnaphthalene	12.12	142	30808	1.695	ng/ul	100
7) Acenaphthylene	14.02	152	48525	1.380	ng/ul#	93
8) Acenaphthene	14.37	153	36968	1.565	ng/ul	99
9) Fluorene	15.36	166	47278	1.595	ng/ul#	97
11) Pentachlorophenol	16.70	266	6711	1.517	ng/ul	98
12) Phenanthrene	17.09	178	79475	1.560	ng/ul	95
13) Anthracene	17.18	178	74746	1.514	ng/ul	95
15) Fluoranthene	19.11	202	102423	1.384	ng/ul	93
17) Pyrene	19.47	202	100000	1.407	ng/ul#	93
18) Benzo(a)anthracene	21.22	228	99681	1.464	ng/ul	97
19) Chrysene	21.27	228	104451	1.540	ng/ul	98
21) Benzo(b)fluoranthene	22.77	252	96496	1.804	ng/ul	89
22) Benzo(k)fluoranthene	22.82	252	101088	1.908	ng/ul#	91
23) Benzo(a)pyrene	23.33	252	81094	1.660	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.61	276	98172	1.604	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.63	278	82670	1.608	ng/ul#	91
26) Benzo(g,h,i)perylene	26.27	276	81627	1.583	ng/ul#	89

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

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