

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM030320\
 Data File : BM025271.D
 Acq On : 05 Mar 2020 04:24
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.434

Quant Time: Mar 05 13:37:00 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 : Wed Mar 04 08:02:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	3437	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	13834	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	9807	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	23646	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	20158	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	16991	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	9232	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	25362	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.49	128	14852	0.357	ng/ul#	82
5) 2-Methylnaphthalene	12.11	142	10740	0.354	ng/ul	99
7) Acenaphthylene	14.02	152	15812	0.344	ng/ul	93
8) Acenaphthene	14.36	153	12691	0.356	ng/ul	99
9) Fluorene	15.35	166	15585	0.349	ng/ul	96
11) Pentachlorophenol	16.70	266	1475	0.238	ng/ul	99
12) Phenanthrene	17.08	178	26952	0.353	ng/ul	96
13) Anthracene	17.17	178	23795	0.352	ng/ul	96
15) Fluoranthene	19.10	202	31888	0.334	ng/ul	94
17) Pyrene	19.46	202	31235	0.409	ng/ul#	94
18) Benzo(a)anthracene	21.21	228	27094	0.366	ng/ul	98
19) Chrysene	21.26	228	29302	0.360	ng/ul	98
21) Benzo(b)fluoranthene	22.76	252	24472	0.360	ng/ul	93
22) Benzo(k)fluoranthene	22.81	252	26826	0.376	ng/ul#	98
23) Benzo(a)pyrene	23.32	252	20332	0.356	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.59	276	25936	0.365	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.61	278	21555	0.366	ng/ul#	94
26) Benzo(g,h,i)perylene	26.26	276	21237	0.352	ng/ul#	91

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

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