

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025298.D
 Acq On : 05 Mar 2020 21:21
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 94 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.436

Quant Time: Mar 06 02:46:04 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	7.67	152	3431	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	14237	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	10085	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	24570	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	22240	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	18199	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	9625	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	27207	0.34	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.49	128	15471	0.361	ng/ul#	82
5) 2-Methylnaphthalene	12.11	142	11184	0.358	ng/ul	99
7) Acenaphthylene	14.01	152	16678	0.353	ng/ul	94
8) Acenaphthene	14.36	153	13285	0.363	ng/ul	99
9) Fluorene	15.35	166	16336	0.356	ng/ul	96
11) Pentachlorophenol	16.70	266	1955	0.304	ng/ul	95
12) Phenanthrene	17.08	178	28344	0.358	ng/ul	95
13) Anthracene	17.17	178	24956	0.355	ng/ul	95
15) Fluoranthene	19.10	202	34043	0.343	ng/ul	94
17) Pyrene	19.46	202	33273	0.394	ng/ul#	94
18) Benzo(a)anthracene	21.21	228	29870	0.365	ng/ul	98
19) Chrysene	21.27	228	32650	0.363	ng/ul	98
21) Benzo(b)fluoranthene	22.77	252	27296	0.375	ng/ul	93
22) Benzo(k)fluoranthene	22.81	252	29293	0.384	ng/ul#	97
23) Benzo(a)pyrene	23.32	252	22152	0.362	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.60	276	26665	0.350	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.61	278	22078	0.350	ng/ul#	95
26) Benzo(g,h,i)perylene	26.26	276	21584	0.334	ng/ul#	92

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

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