

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025313.D
 Acq On : 06 Mar 2020 13:26
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 111 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.438

Quant Time: Mar 06 14:03:20 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	4513	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	18682	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	12836	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	30782	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	29109	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	25839	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	12231	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	33720	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.49	128	19999	0.356	ng/ul#	82
5) 2-Methylnaphthalene	12.11	142	14390	0.351	ng/ul	99
7) Acenaphthylene	14.02	152	20417	0.340	ng/ul#	93
8) Acenaphthene	14.36	153	16697	0.358	ng/ul	100
9) Fluorene	15.35	166	20343	0.348	ng/ul#	95
11) Pentachlorophenol	16.70	266	2662	0.330	ng/ul	98
12) Phenanthrene	17.08	178	35616	0.359	ng/ul	95
13) Anthracene	17.17	178	31049	0.352	ng/ul	96
15) Fluoranthene	19.10	202	42805	0.345	ng/ul	96
17) Pyrene	19.46	202	42735	0.387	ng/ul	97
18) Benzo(a)anthracene	21.22	228	39343	0.368	ng/ul	98
19) Chrysene	21.27	228	42660	0.363	ng/ul	97
21) Benzo(b)fluoranthene	22.77	252	36887	0.357	ng/ul	92
22) Benzo(k)fluoranthene	22.82	252	41565	0.383	ng/ul	99
23) Benzo(a)pyrene	23.33	252	31427	0.362	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.60	276	38838	0.359	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.62	278	32100	0.358	ng/ul#	94
26) Benzo(g,h,i)perylene	26.27	276	30894	0.336	ng/ul#	90

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

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