

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025308.D
 Acq On : 06 Mar 2020 05:46
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 106 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.437

Quant Time: Mar 06 11:33:52 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	3624	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	14818	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	10548	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	24950	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	19676	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	17073	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	9958	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	26158	0.32	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.49	128	15982	0.358	ng/ul#	83
5) 2-Methylnaphthalene	12.11	142	11639	0.358	ng/ul	100
7) Acenaphthylene	14.01	152	16880	0.342	ng/ul	94
8) Acenaphthene	14.36	153	13820	0.361	ng/ul	99
9) Fluorene	15.35	166	16933	0.353	ng/ul	96
11) Pentachlorophenol	16.70	266	2004	0.307	ng/ul	97
12) Phenanthrene	17.08	178	28662	0.356	ng/ul	95
13) Anthracene	17.17	178	25157	0.352	ng/ul	95
15) Fluoranthene	19.10	202	32650	0.324	ng/ul	96
17) Pyrene	19.46	202	31957	0.428	ng/ul#	94
18) Benzo(a)anthracene	21.21	228	26680	0.369	ng/ul	98
19) Chrysene	21.26	228	28802	0.362	ng/ul	98
21) Benzo(b)fluoranthene	22.76	252	24348	0.356	ng/ul	94
22) Benzo(k)fluoranthene	22.81	252	26485	0.370	ng/ul	97
23) Benzo(a)pyrene	23.32	252	20842	0.363	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.60	276	27503	0.385	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.61	278	22672	0.383	ng/ul#	94
26) Benzo(g,h,i)perylene	26.25	276	22564	0.372	ng/ul#	90

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

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