

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022520\
 Data File : BM025082.D
 Acq On : 26 Feb 2020 17:28
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.487

Quant Time: Feb 27 05:08:04 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 25 11:03:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	2329	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	8917	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.32	164	5540	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	12782	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	14114	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	15860	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	6484	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	17277	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.51	128	10926	0.403	ng/ul#	83
5) 2-Methylnaphthalene	12.13	142	7521	0.384	ng/ul	99
7) Acenaphthylene	14.04	152	11781	0.369	ng/ul	94
8) Acenaphthene	14.38	153	8484	0.395	ng/ul	99
9) Fluorene	15.37	166	10143	0.377	ng/ul	96
11) Pentachlorophenol	16.72	266	1437	0.368	ng/ul	98
12) Phenanthrene	17.10	178	18235	0.406	ng/ul	97
13) Anthracene	17.19	178	16222	0.372	ng/ul	96
15) Fluoranthene	19.12	202	22669	0.347	ng/ul	96
17) Pyrene	19.48	202	22751	0.358	ng/ul	98
18) Benzo(a)anthracene	21.23	228	22918	0.376	ng/ul	98
19) Chrysene	21.28	228	24088	0.397	ng/ul	98
21) Benzo(b)fluoranthene	22.80	252	25013	0.402	ng/ul	93
22) Benzo(k)fluoranthene	22.84	252	25185	0.409	ng/ul#	95
23) Benzo(a)pyrene	23.36	252	22508	0.397	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.64	276	28816	0.405	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.66	278	24145	0.404	ng/ul#	95
26) Benzo(g,h,i)perylene	26.31	276	24466	0.408	ng/ul#	92

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

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