

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040320\
 Data File : BM025904.D
 Acq On : 02 Apr 2020 16:28
 Operator : CG/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.435

Quant Time: Apr 02 17:00:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 02 10:49:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	2824	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	11477	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	7145	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	18628	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	18955	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	19081	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	7857	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	23272	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	12208	0.362	ng/ul	99
5) 2-Methylnaphthalene	12.00	142	8988	0.381	ng/ul	100
7) Acenaphthylene	13.92	152	12774	0.418	ng/ul	99
8) Acenaphthene	14.26	153	10085	0.398	ng/ul	99
9) Fluorene	15.26	166	12535	0.399	ng/ul	99
11) Pentachlorophenol	16.61	266	1609	0.352	ng/ul	98
12) Phenanthrene	16.99	178	21949	0.374	ng/ul	99
13) Anthracene	17.08	178	19141	0.369	ng/ul	99
15) Fluoranthene	19.01	202	26719	0.368	ng/ul	99
17) Pyrene	19.38	202	26817	0.395	ng/ul	99
18) Benzo(a)anthracene	21.13	228	25494	0.376	ng/ul	99
19) Chrysene	21.18	228	28358	0.383	ng/ul	100
21) Benzo(b)fluoranthene	22.65	252	27687	0.394	ng/ul	99
22) Benzo(k)fluoranthene	22.70	252	29126	0.387	ng/ul	100
23) Benzo(a)pyrene	23.20	252	25417	0.405	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.41	276	30114	0.358	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.42	278	25020	0.357	ng/ul	100
26) Benzo(g,h,i)perylene	26.06	276	24849	0.350	ng/ul	100

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

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