

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040720\
 Data File : BM025952.D
 Acq On : 07 Apr 2020 10:35
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.442

Quant Time: Apr 07 11:06:37 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	7.56	152	1940	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	8459	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	5222	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	12921	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	12499	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	12512	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	5829	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	15790	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	9119	0.367	ng/ul	99
5) 2-Methylnaphthalene	12.00	142	6640	0.382	ng/ul	98
7) Acenaphthylene	13.92	152	9595	0.430	ng/ul	98
8) Acenaphthene	14.26	153	7378	0.398	ng/ul	99
9) Fluorene	15.26	166	8948	0.389	ng/ul	100
11) Pentachlorophenol	16.61	266	957	0.302	ng/ul	97
12) Phenanthrene	16.99	178	15272	0.375	ng/ul	99
13) Anthracene	17.08	178	13307	0.370	ng/ul	100
15) Fluoranthene	19.01	202	18313	0.364	ng/ul	99
17) Pyrene	19.37	202	18474	0.413	ng/ul	97
18) Benzo(a)anthracene	21.13	228	16957	0.380	ng/ul	100
19) Chrysene	21.18	228	19020	0.390	ng/ul	99
21) Benzo(b)fluoranthene	22.65	252	18321	0.398	ng/ul	97
22) Benzo(k)fluoranthene	22.69	252	19644	0.398	ng/ul	99
23) Benzo(a)pyrene	23.20	252	16966	0.412	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.41	276	18672	0.339	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.42	278	15238	0.331	ng/ul	97
26) Benzo(g,h,i)perylene	26.05	276	15744	0.338	ng/ul	97

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

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