

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM040420\
 Data File : BM025906.D
 Acq On : 03 Apr 2020 08:53
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.436

Quant Time: Apr 03 10:17:03 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	2188	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	8860	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	5545	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	14586	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	14995	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	15475	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	6066	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	18206	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	9549	0.366	ng/ul	100
5) 2-Methylnaphthalene	12.00	142	6950	0.381	ng/ul	98
7) Acenaphthylene	13.92	152	9998	0.422	ng/ul	100
8) Acenaphthene	14.26	153	7956	0.404	ng/ul	99
9) Fluorene	15.26	166	9871	0.404	ng/ul	98
11) Pentachlorophenol	16.61	266	1180	0.330	ng/ul	99
12) Phenanthrene	16.99	178	17271	0.376	ng/ul	99
13) Anthracene	17.08	178	14711	0.362	ng/ul	99
15) Fluoranthene	19.01	202	21061	0.371	ng/ul	100
17) Pyrene	19.37	202	21118	0.393	ng/ul	98
18) Benzo(a)anthracene	21.13	228	19289	0.360	ng/ul	99
19) Chrysene	21.18	228	22712	0.388	ng/ul	100
21) Benzo(b)fluoranthene	22.66	252	22370	0.393	ng/ul	99
22) Benzo(k)fluoranthene	22.70	252	23984	0.393	ng/ul	99
23) Benzo(a)pyrene	23.20	252	20540	0.404	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.41	276	26317	0.386	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.42	278	21791	0.383	ng/ul	99
26) Benzo(g,h,i)perylene	26.05	276	22089	0.383	ng/ul	100

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

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