

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080719\
 Data File : BM021915.D
 Acq On : 08 Aug 2019 02:47
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.441

Quant Time: Aug 08 06:54:45 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM080319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 08 04:39:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	684	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	3115	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	1683	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	3690	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	3455	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	3991	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	1977	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	4388	0.35	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.47	128	3251	0.369	ng/ul	100
5) 2-Methylnaphthalene	12.11	142	2238	0.382	ng/ul	99
7) Acenaphthylene	14.01	152	2973	0.379	ng/ul	99
8) Acenaphthene	14.35	153	2454	0.382	ng/ul	97
9) Fluorene	15.35	166	2596	0.364	ng/ul	98
11) Pentachlorophenol	16.72	266	264	0.388	ng/ul#	83
12) Phenanthrene	17.09	178	3961	0.369	ng/ul	99
13) Anthracene	17.19	178	3481	0.380	ng/ul	98
15) Fluoranthene	19.11	202	4887	0.340	ng/ul	98
17) Pyrene	19.47	202	5156	0.363	ng/ul#	95
18) Benzo(a)anthracene	21.23	228	4549	0.368	ng/ul	99
19) Chrysene	21.28	228	5615	0.353	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	4974	0.370	ng/ul	99
22) Benzo(k)fluoranthene	22.83	252	5456	0.358	ng/ul	98
23) Benzo(a)pyrene	23.36	252	4981	0.364	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.68	276	5948	0.364	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.69	278	4775	0.362	ng/ul	98
26) Benzo(g,h,i)perylene	26.35	276	5248	0.357	ng/ul	96

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

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