

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM080319\
 Data File : BM021814.D
 Acq On : 04 Aug 2019 16:00
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.428

Quant Time: Aug 05 05:30:42 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 : Mon Aug 05 05:26:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	646	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	2813	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	1713	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	4129	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	4080	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	4829	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	1967	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	5470	0.39	ng/ul	0.00

Target Compounds						Qvalue
3) Naphthalene	10.48	128	3225	0.406	ng/ul	98
5) 2-Methylnaphthalene	12.11	142	2266	0.428	ng/ul	93
7) Acenaphthylene	14.02	152	3329	0.417	ng/ul	98
8) Acenaphthene	14.36	153	2707	0.414	ng/ul	96
9) Fluorene	15.36	166	2962	0.408	ng/ul	98
11) Pentachlorophenol	16.73	266	333	0.437	ng/ul	96
12) Phenanthrene	17.10	178	4788	0.399	ng/ul	99
13) Anthracene	17.20	178	4231	0.413	ng/ul	98
15) Fluoranthene	19.12	202	6301	0.392	ng/ul	100
17) Pyrene	19.48	202	6689	0.399	ng/ul	99
18) Benzo(a)anthracene	21.23	228	6040	0.414	ng/ul	99
19) Chrysene	21.28	228	7363	0.392	ng/ul	99
21) Benzo(b)fluoranthene	22.80	252	6632	0.407	ng/ul	97
22) Benzo(k)fluoranthene	22.85	252	7353	0.399	ng/ul	99
23) Benzo(a)pyrene	23.37	252	6690	0.404	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.69	276	8104	0.410	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.70	278	6492	0.406	ng/ul	98
26) Benzo(g,h,i)perylene	26.37	276	7160	0.402	ng/ul	99

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

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