

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080319\
 Data File : BM021856.D
 Acq On : 05 Aug 2019 23:33
 Operator : HP/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.433

Quant Time: Aug 06 01:32:51 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 : Tue Aug 06 01:29:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	846	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	3652	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	2023	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	4415	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	4427	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	5141	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	2506	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	5929	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	4119	0.399	ng/ul	96
5) 2-Methylnaphthalene	12.10	142	2884	0.420	ng/ul	98
7) Acenaphthylene	14.01	152	3887	0.412	ng/ul	97
8) Acenaphthene	14.35	153	3188	0.413	ng/ul	97
9) Fluorene	15.35	166	3422	0.399	ng/ul	98
11) Pentachlorophenol	16.72	266	362	0.444	ng/ul	94
12) Phenanthrene	17.09	178	5318	0.414	ng/ul	99
13) Anthracene	17.18	178	4815	0.439	ng/ul	97
15) Fluoranthene	19.11	202	6818	0.396	ng/ul	100
17) Pyrene	19.48	202	7317	0.402	ng/ul	97
18) Benzo(a)anthracene	21.23	228	6737	0.425	ng/ul	99
19) Chrysene	21.28	228	7826	0.384	ng/ul	98
21) Benzo(b)fluoranthene	22.79	252	7269	0.419	ng/ul	96
22) Benzo(k)fluoranthene	22.84	252	7605	0.388	ng/ul	98
23) Benzo(a)pyrene	23.36	252	7165	0.406	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.68	276	8799	0.418	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.69	278	7056	0.415	ng/ul	97
26) Benzo(g,h,i)perylene	26.36	276	7448	0.393	ng/ul	100

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

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