

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080619\
 Data File : BM021862.D
 Acq On : 06 Aug 2019 17:21
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.435

Quant Time: Aug 06 17:56:39 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	714	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	3196	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	1799	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	3962	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	3698	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	4362	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	2208	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	5069	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	3606	0.399	ng/ul	98
5) 2-Methylnaphthalene	12.11	142	2516	0.418	ng/ul	99
7) Acenaphthylene	14.02	152	3404	0.406	ng/ul	98
8) Acenaphthene	14.36	153	2793	0.407	ng/ul	98
9) Fluorene	15.35	166	3003	0.394	ng/ul	98
11) Pentachlorophenol	16.72	266	322	0.440	ng/ul	84
12) Phenanthrene	17.09	178	4678	0.406	ng/ul	99
13) Anthracene	17.19	178	4153	0.422	ng/ul	99
15) Fluoranthene	19.12	202	5807	0.376	ng/ul	99
17) Pyrene	19.48	202	6141	0.404	ng/ul	97
18) Benzo(a)anthracene	21.23	228	5394	0.408	ng/ul	99
19) Chrysene	21.28	228	6706	0.394	ng/ul	100
21) Benzo(b)fluoranthene	22.80	252	6015	0.409	ng/ul	96
22) Benzo(k)fluoranthene	22.84	252	6504	0.391	ng/ul	97
23) Benzo(a)pyrene	23.37	252	6002	0.401	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	25.69	276	7089	0.397	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.70	278	5745	0.398	ng/ul	96
26) Benzo(g,h,i)perylene	26.37	276	6156	0.383	ng/ul	98

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

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