

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080219\
 Data File : BM021763.D
 Acq On : 02 Aug 2019 18:33
 Operator : HP/JU
 Sample : SST0.415
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.415

Quant Time: Aug 02 23:41:12 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM080219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 02 23:36:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	1076	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	4325	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	1711	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.08	188	3262	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	2433	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	2170	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	2307	0.45	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	2873	0.31	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.51	128	4780	0.438	ng/ul	100
5) 2-Methylnaphthalene	12.14	142	2767	0.480	ng/ul	100
7) Acenaphthylene	14.03	152	3374	0.361	ng/ul	100
8) Acenaphthene	14.38	153	2868	0.416	ng/ul	100
9) Fluorene	15.38	166	3004	0.388	ng/ul	100
11) Pentachlorophenol	16.76	266	175	0.157	ng/ul	98
12) Phenanthrene	17.12	178	4302	0.477	ng/ul	100
13) Anthracene	17.22	178	3244	0.418	ng/ul	100
15) Fluoranthene	19.13	202	4057	0.360	ng/ul	100
17) Pyrene	19.50	202	4242	0.616	ng/ul	100
18) Benzo(a)anthracene	21.25	228	2717	0.358	ng/ul	100
19) Chrysene	21.30	228	4471	0.468	ng/ul	100
21) Benzo(b)fluoranthene	22.82	252	3563	0.445	ng/ul	100
22) Benzo(k)fluoranthene	22.87	252	3962	0.488	ng/ul	100
23) Benzo(a)pyrene	23.40	252	2925	0.406	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.73	276	3542	0.368	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.75	278	2784	0.369	ng/ul	100
26) Benzo(g,h,i)perylene	26.41	276	3434	0.372	ng/ul	100

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

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