

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM080319\
 Data File : BM021781.D
 Acq On : 03 Aug 2019 18:42
 Operator : HP/JU
 Sample : SSTD0.421
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.421

Quant Time: Aug 03 22:38:18 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 : Sat Aug 03 22:35:19 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	712	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	2889	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	1612	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	3790	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	3811	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	4349	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	1875	0.47	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	5221	0.54	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.49	128	3257	0.401	ng/ul	100
5) 2-Methylnaphthalene	12.12	142	2167	0.451	ng/ul	100
7) Acenaphthylene	14.02	152	3078	0.383	ng/ul	100
8) Acenaphthene	14.36	153	2474	0.366	ng/ul	100
9) Fluorene	15.37	166	2780	0.388	ng/ul	100
11) Pentachlorophenol	16.74	266	254	0.457	ng/ul	100
12) Phenanthrene	17.10	178	4454	0.359	ng/ul	100
13) Anthracene	17.20	178	3767	0.390	ng/ul	100
15) Fluoranthene	19.12	202	5932	0.448	ng/ul	100
17) Pyrene	19.49	202	6379	0.381	ng/ul	100
18) Benzo(a)anthracene	21.24	228	5521	0.459	ng/ul	100
19) Chrysene	21.29	228	7017	0.366	ng/ul	100
21) Benzo(b)fluoranthene	22.81	252	6040	0.321	ng/ul	100
22) Benzo(k)fluoranthene	22.85	252	6754	0.340	ng/ul	100
23) Benzo(a)pyrene	23.38	252	5994	0.392	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.70	276	7201	0.408	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.72	278	5811	0.419	ng/ul	100
26) Benzo(g,h,i)perylene	26.39	276	6578	0.392	ng/ul	100

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

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