

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022420\
 Data File : BM025030.D
 Acq On : 24 Feb 2020 20:11
 Operator : CG/JU
 Sample : SST0.479
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.479

Quant Time: Feb 25 00:51:38 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 25 00:46:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	1720	0.40	ng/ul	0.00
2) Naphthalene-d8	10.47	136	6678	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	4560	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	11184	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	12284	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	13381	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	5078	0.46	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	15437	0.50	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.52	128	8172	0.429	ng/ul	100
5) 2-Methylnaphthalene	12.14	142	5943	0.456	ng/ul	100
7) Acenaphthylene	14.05	152	10467	0.496	ng/ul	100
8) Acenaphthene	14.39	153	7091	0.401	ng/ul	100
9) Fluorene	15.38	166	8875	0.438	ng/ul	100
11) Pentachlorophenol	16.73	266	1338	0.590	ng/ul	100
12) Phenanthrene	17.11	178	15333	0.430	ng/ul	100
13) Anthracene	17.20	178	15152	0.496	ng/ul	100
15) Fluoranthene	19.13	202	21413	0.557	ng/ul	100
17) Pyrene	19.49	202	21464	0.385	ng/ul	100
18) Benzo(a)anthracene	21.24	228	20945	0.446	ng/ul	100
19) Chrysene	21.29	228	20895	0.422	ng/ul	100
21) Benzo(b)fluoranthene	22.80	252	21040	0.362	ng/ul	100
22) Benzo(k)fluoranthene	22.84	252	20974	0.354	ng/ul	100
23) Benzo(a)pyrene	23.36	252	19231	0.378	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.65	276	24309	0.392	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.67	278	20393	0.407	ng/ul	100
26) Benzo(g,h,i)perylene	26.32	276	20334	0.411	ng/ul	100

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

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