

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040521\
 Data File : BM029340.D
 Acq On : 05 Apr 2021 13:22
 Operator : CG/JU
 Sample : SST0.478
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.478

Quant Time: Apr 05 14:10:05 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 05 14:08:43 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	2079	0.40	ng/ul	0.00
2) Naphthalene-d8	10.48	136	6781	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	3516	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	7311	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	7858	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	8313	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	3837	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	7699	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.53	128	7702	0.456	ng/ul	100
5) 2-Methylnaphthalene	12.15	142	4952	0.438	ng/ul	100
7) Acenaphthylene	14.05	152	6111	0.426	ng/ul	100
8) Acenaphthene	14.39	153	5202	0.457	ng/ul	100
9) Fluorene	15.38	166	5732	0.454	ng/ul	100
11) Pentachlorophenol	16.73	266	681	0.356	ng/ul	99
12) Phenanthrene	17.11	178	8980	0.439	ng/ul	100
13) Anthracene	17.20	178	7936	0.407	ng/ul	100
15) Fluoranthene	19.13	202	9659	0.415	ng/ul	100
17) Pyrene	19.49	202	10060	0.352	ng/ul	100
18) Benzo(a)anthracene	21.23	228	8854	0.330	ng/ul	100
19) Chrysene	21.28	228	9596	0.360	ng/ul	100
21) Benzo(b)fluoranthene	22.79	252	10066	0.354	ng/ul	100
22) Benzo(k)fluoranthene	22.84	252	10528	0.371	ng/ul	100
23) Benzo(a)pyrene	23.36	252	8810	0.348	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.69	276	12207	0.370	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.70	278	10291	0.372	ng/ul	100
26) Benzo(g,h,i)perylene	26.36	276	10310	0.375	ng/ul	100

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

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