

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041520\
 Data File : BM025958.D
 Acq On : 15 Apr 2020 10:03
 Operator : MA/SJ
 Sample : SST0.476
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.476

Quant Time: Apr 15 10:43:19 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 15 10:41:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	1121	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	4233	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	2667	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	6133	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	5171	0.40	ng/ul	0.00
20) Perylene-d12	23.28	264	4603	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	2353	0.31	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	5688	0.30	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	4435	0.356	ng/ul	100
5) 2-Methylnaphthalene	12.00	142	2989	0.343	ng/ul	100
7) Acenaphthylene	13.92	152	3966	0.348	ng/ul	100
8) Acenaphthene	14.26	153	3419	0.361	ng/ul	100
9) Fluorene	15.25	166	4026	0.343	ng/ul	100
11) Pentachlorophenol	16.61	266	455	0.302	ng/ul	100
12) Phenanthrene	16.98	178	6775	0.351	ng/ul	100
13) Anthracene	17.08	178	5734	0.336	ng/ul	100
15) Fluoranthene	19.01	202	7842	0.328	ng/ul	100
17) Pyrene	19.37	202	7855	0.424	ng/ul	100
18) Benzo(a)anthracene	21.12	228	6310	0.341	ng/ul	100
19) Chrysene	21.17	228	7892	0.391	ng/ul	100
21) Benzo(b)fluoranthene	22.65	252	6662	0.393	ng/ul	100
22) Benzo(k)fluoranthene	22.69	252	7405	0.408	ng/ul	100
23) Benzo(a)pyrene	23.19	252	5648	0.373	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.40	276	6272	0.309	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.42	278	4749	0.281	ng/ul	100
26) Benzo(g,h,i)perylene	26.04	276	5841	0.341	ng/ul	100

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

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