

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040220\
 Data File : BM025888.D
 Acq On : 01 Apr 2020 15:33
 Operator : CG/JU
 Sample : SST0.477
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.477

Quant Time: Apr 01 16:06:13 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 01 16:04:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	2681	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	10564	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	7035	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	17137	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	18352	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	20057	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	6256	0.35	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	17272	0.35	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.38	128	10436	0.332	ng/ul	100
5) 2-Methylnaphthalene	12.00	142	7276	0.344	ng/ul	100
7) Acenaphthylene	13.92	152	9649	0.323	ng/ul	100
8) Acenaphthene	14.27	153	8188	0.328	ng/ul	100
9) Fluorene	15.26	166	10063	0.341	ng/ul	100
11) Pentachlorophenol	16.61	266	1287	0.363	ng/ul	100
12) Phenanthrene	16.99	178	17535	0.326	ng/ul	100
13) Anthracene	17.08	178	15383	0.342	ng/ul	100
15) Fluoranthene	19.01	202	21429	0.344	ng/ul	100
17) Pyrene	19.38	202	21383	0.302	ng/ul	100
18) Benzo(a)anthracene	21.13	228	21189	0.351	ng/ul	100
19) Chrysene	21.18	228	23366	0.321	ng/ul	100
21) Benzo(b)fluoranthene	22.66	252	23971	0.309	ng/ul	100
22) Benzo(k)fluoranthene	22.70	252	25464	0.320	ng/ul	100
23) Benzo(a)pyrene	23.20	252	21125	0.331	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.41	276	28766	0.354	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.43	278	23914	0.381	ng/ul	100
26) Benzo(g,h,i)perylene	26.06	276	24213	0.306	ng/ul	100

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

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