

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025411.D
 Acq On : 09 Mar 2020 11:10
 Operator : CG/JU
 Sample : SST0.481
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.481

Quant Time: Mar 09 12:05:36 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 09 12:03:47 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	2532	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	10006	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	7012	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	16678	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	12728	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	10883	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	6171	0.33	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	15649	0.29	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	10129	0.336	ng/ul	100
5) 2-Methylnaphthalene	12.10	142	7322	0.334	ng/ul	100
7) Acenaphthylene	14.01	152	9603	0.292	ng/ul	100
8) Acenaphthene	14.35	153	8284	0.325	ng/ul	100
9) Fluorene	15.34	166	10266	0.322	ng/ul	100
11) Pentachlorophenol	16.69	266	1107	0.254	ng/ul	100
12) Phenanthrene	17.08	178	17335	0.322	ng/ul	100
13) Anthracene	17.17	178	15000	0.314	ng/ul	100
15) Fluoranthene	19.09	202	19605	0.291	ng/ul	100
17) Pyrene	19.45	202	18990	0.393	ng/ul	100
18) Benzo(a)anthracene	21.21	228	14957	0.320	ng/ul	100
19) Chrysene	21.26	228	16795	0.327	ng/ul	100
21) Benzo(b)fluoranthene	22.75	252	14108	0.324	ng/ul	100
22) Benzo(k)fluoranthene	22.80	252	15012	0.329	ng/ul	100
23) Benzo(a)pyrene	23.31	252	11682	0.320	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.58	276	15522	0.341	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.59	278	12913	0.342	ng/ul	100
26) Benzo(g,h,i)perylene	26.24	276	13381	0.346	ng/ul	100

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

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