

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN060220\
 Data File : BN010743.D
 Acq On : 02 Jun 2020 13:59
 Operator : CG/JU
 Sample : SST00.401
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.401

Quant Time: Jun 02 14:32:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN060220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 02 14:30:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	1918	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	6340	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	3272	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	6587	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	5688	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	5063	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	4040	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	7422	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	7212	0.358	ng/ul	100
5) 2-Methylnaphthalene	11.99	142	4793	0.356	ng/ul	100
7) Acenaphthylene	13.90	152	5729	0.329	ng/ul	100
8) Acenaphthene	14.26	153	4750	0.355	ng/ul	100
9) Fluorene	15.25	166	5309	0.334	ng/ul	100
11) Pentachlorophenol	16.60	266	643	0.500	ng/ul	100
12) Phenanthrene	16.98	178	8383	0.354	ng/ul	100
13) Anthracene	17.07	178	6944	0.333	ng/ul	100
15) Fluoranthene	19.00	202	9146	0.317	ng/ul	100
17) Pyrene	19.37	202	9180	0.340	ng/ul	100
18) Benzo(a)anthracene	21.13	228	8060	0.329	ng/ul	100
19) Chrysene	21.18	228	9672	0.346	ng/ul	100
21) Benzo(b)fluoranthene	22.66	252	8977	0.392	ng/ul	100
22) Benzo(k)fluoranthene	22.71	252	8971	0.371	ng/ul	100
23) Benzo(a)pyrene	23.21	252	7951	0.400	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.42	276	10088	0.405	ng/ul	100
25) Dibenzo(a,h)anthracene	25.44	278	8306	0.408	ng/ul	100
26) Benzo(g,h,i)perylene	26.06	276	9194	0.420	ng/ul	100

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

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