

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060320\
 Data File : BN010751.D
 Acq On : 03 Jun 2020 14:21
 Operator : CG/JU
 Sample : SST0.402
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.402

Quant Time: Jun 03 14:57:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN060320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 03 14:56:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1563	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	5144	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	2683	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	5223	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	4356	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	4052	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	5092	0.61	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	9414	0.62	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	6374	0.402	ng/ul	100
5) 2-Methylnaphthalene	11.99	142	3976	0.390	ng/ul	100
7) Acenaphthylene	13.91	152	4958	0.385	ng/ul	100
8) Acenaphthene	14.26	153	4150	0.393	ng/ul	100
9) Fluorene	15.25	166	4562	0.391	ng/ul	100
11) Pentachlorophenol	16.61	266	483	0.347	ng/ul	100
12) Phenanthrene	16.98	178	7130	0.397	ng/ul	100
13) Anthracene	17.08	178	5821	0.390	ng/ul	100
15) Fluoranthene	19.00	202	7954	0.401	ng/ul	100
17) Pyrene	19.37	202	7938	0.418	ng/ul	100
18) Benzo(a)anthracene	21.13	228	6505	0.394	ng/ul	100
19) Chrysene	21.18	228	8341	0.418	ng/ul	100
21) Benzo(b)fluoranthene	22.66	252	7496	0.408	ng/ul	100
22) Benzo(k)fluoranthene	22.71	252	8271	0.421	ng/ul	100
23) Benzo(a)pyrene	23.21	252	6352	0.402	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.42	276	8764	0.404	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.44	278	7245	0.410	ng/ul	100
26) Benzo(g,h,i)perylene	26.06	276	8122	0.424	ng/ul	100

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

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