

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN060220\
 Data File : BN010744.D
 Acq On : 02 Jun 2020 14:29
 Operator : CG/JU
 Sample : SST0.802
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.802

Quant Time: Jun 02 15:21:28 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	7.57	152	1844	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	6049	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	3134	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	6307	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	5589	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	5127	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	7962	0.78	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	14740	0.75	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.37	128	15268	0.795	ng/ul	99
5) 2-Methylnaphthalene	11.99	142	9766	0.759	ng/ul	100
7) Acenaphthylene	13.90	152	12448	0.747	ng/ul	99
8) Acenaphthene	14.25	153	10031	0.783	ng/ul	96
9) Fluorene	15.25	166	11210	0.737	ng/ul	100
11) Pentachlorophenol	16.60	266	1324	1.076	ng/ul	99
12) Phenanthrene	16.98	178	17681	0.779	ng/ul	98
13) Anthracene	17.07	178	14955	0.749	ng/ul	99
15) Fluoranthene	19.00	202	19838	0.718	ng/ul	100
17) Pyrene	19.37	202	19581	0.738	ng/ul	100
18) Benzo(a)anthracene	21.13	228	16961	0.705	ng/ul	99
19) Chrysene	21.18	228	20392	0.743	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	18881	0.813	ng/ul	94
22) Benzo(k)fluoranthene	22.71	252	20633	0.842	ng/ul	94
23) Benzo(a)pyrene	23.21	252	16110	0.801	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.42	276	22770	0.904	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.43	278	18716	0.907	ng/ul	96
26) Benzo(g,h,i)perylene	26.06	276	19949	0.901	ng/ul	97

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

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