

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN012721\
 Data File : BN013492.D
 Acq On : 26 Jan 2021 17:03
 Operator : CG/JU
 Sample : SST0.804
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.804

Quant Time: Jan 26 17:46:56 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN012721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 26 17:05:29 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	4084	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	14715	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	7568	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	15471	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	13873	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	12747	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	18162	0.83	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	34540	0.82	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	34648	0.814	ng/ul	100
5) 2-Methylnaphthalene	12.00	142	22164	0.792	ng/ul	99
7) Acenaphthylene	13.92	152	25391	0.709	ng/ul	99
8) Acenaphthene	14.26	153	22549	0.767	ng/ul	99
9) Fluorene	15.26	166	25064	0.743	ng/ul	97
11) Pentachlorophenol	16.60	266	1772	0.395	ng/ul	98
12) Phenanthrene	16.99	178	41412	0.798	ng/ul	99
13) Anthracene	17.08	178	34542	0.782	ng/ul	100
15) Fluoranthene	19.02	202	46884	0.761	ng/ul	99
17) Pyrene	19.38	202	46033	0.702	ng/ul	100
18) Benzo(a)anthracene	21.14	228	37803	0.659	ng/ul	99
19) Chrysene	21.19	228	45926	0.724	ng/ul	100
21) Benzo(b)fluoranthene	22.68	252	43711	0.741	ng/ul	97
22) Benzo(k)fluoranthene	22.72	252	45912	0.704	ng/ul	97
23) Benzo(a)pyrene	23.23	252	36847	0.673	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.48	276	48817	0.679	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.50	278	39903	0.702	ng/ul	98
26) Benzo(g,h,i)perylene	26.13	276	42603	0.655	ng/ul	100

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

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