

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080719\
 Data File : BN006981.D
 Acq On : 07 Aug 2019 11:18
 Operator : HP/JU
 Sample : SST00.221
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.221

Quant Time: Aug 07 17:25:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN080719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 17:13:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	4933	0.40	ng/ul	0.00
2) Naphthalene-d8	10.24	136	18468	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.12	164	11567	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	29583	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	27315	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	28398	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.84	152	6667	0.20	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	19631	0.21	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.28	128	10764	0.203	ng/ul	97
5) 2-Methylnaphthalene	11.91	142	8060	0.202	ng/ul	99
7) Acenaphthylene	13.83	152	13577	0.203	ng/ul	99
8) Acenaphthene	14.18	153	9353	0.204	ng/ul	98
9) Fluorene	15.17	166	11417	0.205	ng/ul	99
11) Pentachlorophenol	16.52	266	1819	0.176	ng/ul	93
12) Phenanthrene	16.90	178	19178	0.203	ng/ul	99
13) Anthracene	17.00	178	19659	0.204	ng/ul	100
15) Fluoranthene	18.94	202	24133	0.206	ng/ul	100
17) Pyrene	19.30	202	24703	0.203	ng/ul	99
18) Benzo(a)anthracene	21.07	228	24310	0.203	ng/ul	100
19) Chrysene	21.12	228	22968	0.204	ng/ul	100
21) Benzo(b)fluoranthene	22.59	252	23802	0.207	ng/ul	97
22) Benzo(k)fluoranthene	22.63	252	22499	0.202	ng/ul#	97
23) Benzo(a)pyrene	23.12	252	22001	0.203	ng/ul#	96
24) Indeno(1,2,3-cd)pyrene	25.32	276	23123	0.198	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.33	278	18379	0.197	ng/ul	95
26) Benzo(g,h,i)perylene	25.95	276	19354	0.197	ng/ul	97

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

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