

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN020421\
 Data File : BN013569.D
 Acq On : 04 Feb 2021 13:43
 Operator : CG/JU
 Sample : SSTD1.606
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.606

Quant Time: Feb 04 14:12:34 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN020421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 04 13:01:49 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	4677	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	16983	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	8993	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	18220	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	17258	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	13112	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	38152	1.50	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	72452	1.50	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.36	128	68479	1.394	ng/ul	99
5) 2-Methylnaphthalene	11.98	142	46015	1.475	ng/ul	100
7) Acenaphthylene	13.90	152	53166	1.435	ng/ul	100
8) Acenaphthene	14.25	153	46105	1.398	ng/ul	99
9) Fluorene	15.25	166	52336	1.435	ng/ul	99
11) Pentachlorophenol	16.59	266	3715	1.383	ng/ul	100
12) Phenanthrene	16.98	178	84353	1.414	ng/ul	99
13) Anthracene	17.07	178	72775	1.487	ng/ul	99
15) Fluoranthene	19.01	202	93308	1.431	ng/ul	99
17) Pyrene	19.37	202	91277	1.292	ng/ul	100
18) Benzo(a)anthracene	21.13	228	87562	1.537	ng/ul	99
19) Chrysene	21.18	228	92354	1.300	ng/ul	100
21) Benzo(b)fluoranthene	22.66	252	85715	1.557	ng/ul	97
22) Benzo(k)fluoranthene	22.71	252	87314	1.489	ng/ul	96
23) Benzo(a)pyrene	23.21	252	68694	1.471	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.46	276	92737	1.495	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.47	278	77069	1.538	ng/ul	97
26) Benzo(g,h,i)perylene	26.11	276	79802	1.453	ng/ul	99

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

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