

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022121\
 Data File : BN013699.D
 Acq On : 20 Feb 2021 21:38
 Operator : CG/JU
 Sample : SSTD3.202
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD3.202

Quant Time: Feb 20 23:13:44 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN022121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 20 23:07:21 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	4361	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	16985	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	9581	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	20460	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	19617	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	18216	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	76524	3.16	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	172025	3.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	133043	3.009	ng/ul	99
5) 2-Methylnaphthalene	11.98	142	91478	3.087	ng/ul	99
7) Acenaphthylene	13.90	152	119542	3.347	ng/ul	99
8) Acenaphthene	14.25	153	96680	3.043	ng/ul	99
9) Fluorene	15.24	166	112348	3.125	ng/ul	100
11) Pentachlorophenol	16.59	266	15696	6.301	ng/ul	97
12) Phenanthrene	16.97	178	182125	2.990	ng/ul	99
13) Anthracene	17.06	178	170135	3.369	ng/ul	99
15) Fluoranthene	19.00	202	210257	3.202	ng/ul	100
17) Pyrene	19.37	202	212686	3.016	ng/ul	97
18) Benzo(a)anthracene	21.12	228	203740	3.200	ng/ul	99
19) Chrysene	21.17	228	202299	2.920	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	210969	2.833	ng/ul	90
22) Benzo(k)fluoranthene	22.70	252	209456	2.715	ng/ul#	91
23) Benzo(a)pyrene	23.21	252	184361	3.080	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.45	276	242817	3.049	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.47	278	197972	3.029	ng/ul	94
26) Benzo(g,h,i)perylene	26.11	276	206343	2.949	ng/ul	98

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

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