

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022121\
 Data File : BN013701.D
 Acq On : 20 Feb 2021 23:29
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.417

Quant Time: Feb 21 00:00:11 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 23:16:49 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	3495	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	12882	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	7363	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	15999	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	15173	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	14025	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	8150	0.43	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	18181	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.35	128	15043	0.443	ng/ul	99
5) 2-Methylnaphthalene	11.98	142	9951	0.431	ng/ul	100
7) Acenaphthylene	13.90	152	12200	0.426	ng/ul	99
8) Acenaphthene	14.25	153	10700	0.433	ng/ul	99
9) Fluorene	15.24	166	12277	0.428	ng/ul	99
11) Pentachlorophenol	16.59	266	1408	0.378	ng/ul	97
12) Phenanthrene	16.97	178	20709	0.431	ng/ul	100
13) Anthracene	17.06	178	17886	0.424	ng/ul	99
15) Fluoranthene	19.00	202	23699	0.432	ng/ul	99
17) Pyrene	19.37	202	24375	0.420	ng/ul	98
18) Benzo(a)anthracene	21.12	228	22411	0.429	ng/ul	99
19) Chrysene	21.17	228	23758	0.440	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	23875	0.445	ng/ul	98
22) Benzo(k)fluoranthene	22.70	252	23841	0.432	ng/ul	99
23) Benzo(a)pyrene	23.21	252	21706	0.457	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.45	276	26980	0.441	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.46	278	21988	0.440	ng/ul	99
26) Benzo(g,h,i)perylene	26.10	276	23361	0.442	ng/ul	100

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

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