

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022021\
 Data File : BN013692.D
 Acq On : 20 Feb 2021 15:27
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.416

Quant Time: Feb 20 16:31:07 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 : Tue Feb 16 16:31:21 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	3927	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	15672	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	8765	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	18597	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	16450	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	15846	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	9886	0.44	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	21712	0.48	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.36	128	18108	0.444	ng/ul	99
5) 2-Methylnaphthalene	11.98	142	11969	0.438	ng/ul	100
7) Acenaphthylene	13.90	152	15862	0.485	ng/ul	99
8) Acenaphthene	14.25	153	12927	0.445	ng/ul	99
9) Fluorene	15.24	166	14653	0.446	ng/ul	99
11) Pentachlorophenol	16.59	266	2429	1.073	ng/ul	99
12) Phenanthrene	16.98	178	23718	0.428	ng/ul	99
13) Anthracene	17.07	178	22888	0.499	ng/ul	99
15) Fluoranthene	19.00	202	27866	0.467	ng/ul	98
17) Pyrene	19.37	202	28512	0.482	ng/ul	99
18) Benzo(a)anthracene	21.12	228	25529	0.478	ng/ul	95
19) Chrysene	21.18	228	25348	0.436	ng/ul	94
21) Benzo(b)fluoranthene	22.66	252	26111	0.403	ng/ul	83
22) Benzo(k)fluoranthene	22.71	252	25444	0.379	ng/ul#	84
23) Benzo(a)pyrene	23.21	252	23614	0.454	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	25.46	276	29286	0.423	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.47	278	23987	0.422	ng/ul#	87
26) Benzo(g,h,i)perylene	26.12	276	25079	0.412	ng/ul	92

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

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