

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN021621\
 Data File : BN013655.D
 Acq On : 17 Feb 2021 00:44
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.414

Quant Time: Feb 17 05:15:17 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	5128	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	19548	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	10312	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	21086	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	18602	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	17158	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	11638	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	21742	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	21368	0.420	ng/ul	99
5) 2-Methylnaphthalene	11.98	142	14124	0.414	ng/ul	100
7) Acenaphthylene	13.90	152	15940	0.415	ng/ul	100
8) Acenaphthene	14.25	153	14265	0.417	ng/ul	99
9) Fluorene	15.24	166	15677	0.405	ng/ul	99
11) Pentachlorophenol	16.59	266	1291	0.503	ng/ul	99
12) Phenanthrene	16.98	178	25669	0.409	ng/ul	100
13) Anthracene	17.07	178	21466	0.413	ng/ul	99
15) Fluoranthene	19.00	202	27247	0.403	ng/ul	97
17) Pyrene	19.37	202	27589	0.413	ng/ul	99
18) Benzo(a)anthracene	21.12	228	23765	0.394	ng/ul	99
19) Chrysene	21.17	228	27674	0.421	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	27188	0.388	ng/ul	98
22) Benzo(k)fluoranthene	22.70	252	29092	0.400	ng/ul	99
23) Benzo(a)pyrene	23.21	252	23635	0.419	ng/ul#	96
24) Indeno(1,2,3-cd)pyrene	25.45	276	30866	0.411	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.46	278	24998	0.406	ng/ul	99
26) Benzo(g,h,i)perylene	26.10	276	26170	0.397	ng/ul	100

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

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