

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012721\
 Data File : BN013499.D
 Acq On : 27 Jan 2021 09:59
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.408

Quant Time: Jan 27 11:56:18 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN012721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 27 11:55:49 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	4449	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	15966	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	8022	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	16429	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	14643	0.40	ng/ul	0.00
20) Perylene-d12	23.32	264	14820	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	9979	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	18104	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	19311	0.418	ng/ul	99
5) 2-Methylnaphthalene	12.00	142	12238	0.417	ng/ul	100
7) Acenaphthylene	13.92	152	13731	0.415	ng/ul	100
8) Acenaphthene	14.26	153	12282	0.417	ng/ul	99
9) Fluorene	15.25	166	13520	0.416	ng/ul	100
11) Pentachlorophenol	16.60	266	847	0.350	ng/ul	96
12) Phenanthrene	16.99	178	22376	0.416	ng/ul	99
13) Anthracene	17.08	178	18338	0.416	ng/ul	100
15) Fluoranthene	19.02	202	24151	0.411	ng/ul	98
17) Pyrene	19.38	202	24433	0.408	ng/ul	99
18) Benzo(a)anthracene	21.14	228	19837	0.410	ng/ul	99
19) Chrysene	21.19	228	25000	0.415	ng/ul	99
21) Benzo(b)fluoranthene	22.67	252	23777	0.382	ng/ul	99
22) Benzo(k)fluoranthene	22.71	252	27518	0.415	ng/ul	99
23) Benzo(a)pyrene	23.23	252	21597	0.409	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.48	276	29971	0.427	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.49	278	24043	0.424	ng/ul	98
26) Benzo(g,h,i)perylene	26.13	276	25157	0.405	ng/ul	99

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

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