

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN021621\
 Data File : BN013640.D
 Acq On : 16 Feb 2021 15:47
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.413

Quant Time: Feb 16 16:16:27 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 10:14:45 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	5034	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	18985	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	9733	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	19258	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	16036	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	14414	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	11222	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	19263	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	20732	0.419	ng/ul	99
5) 2-Methylnaphthalene	11.98	142	13479	0.407	ng/ul	100
7) Acenaphthylene	13.90	152	14843	0.409	ng/ul	99
8) Acenaphthene	14.25	153	13339	0.413	ng/ul	99
9) Fluorene	15.24	166	14515	0.397	ng/ul	98
11) Pentachlorophenol	16.59	266	1071	0.457	ng/ul	100
12) Phenanthrene	16.98	178	23221	0.405	ng/ul	99
13) Anthracene	17.07	178	19255	0.405	ng/ul	99
15) Fluoranthene	19.00	202	24011	0.388	ng/ul	100
17) Pyrene	19.37	202	24254	0.421	ng/ul	99
18) Benzo(a)anthracene	21.12	228	19724	0.379	ng/ul	100
19) Chrysene	21.17	228	24142	0.426	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	22763	0.386	ng/ul	98
22) Benzo(k)fluoranthene	22.70	252	25094	0.411	ng/ul	100
23) Benzo(a)pyrene	23.21	252	19467	0.411	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.45	276	25973	0.412	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.47	278	20920	0.405	ng/ul	99
26) Benzo(g,h,i)perylene	26.10	276	22051	0.398	ng/ul	99

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

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