

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012721\
 Data File : BN013510.D
 Acq On : 27 Jan 2021 17:45
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.409

Quant Time: Jan 27 18:19:07 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	4495	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	16222	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	8401	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	18265	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	15735	0.40	ng/ul	0.00
20) Perylene-d12	23.32	264	13617	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	10001	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	19380	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	19104	0.407	ng/ul	100
5) 2-Methylnaphthalene	12.00	142	12210	0.410	ng/ul	100
7) Acenaphthylene	13.92	152	14299	0.413	ng/ul	99
8) Acenaphthene	14.26	153	12480	0.405	ng/ul	99
9) Fluorene	15.25	166	14488	0.425	ng/ul	99
11) Pentachlorophenol	16.60	266	1521	0.565	ng/ul	99
12) Phenanthrene	16.99	178	24239	0.405	ng/ul	99
13) Anthracene	17.08	178	20256	0.413	ng/ul	99
15) Fluoranthene	19.01	202	25801	0.395	ng/ul	99
17) Pyrene	19.38	202	26016	0.404	ng/ul	99
18) Benzo(a)anthracene	21.13	228	21219	0.409	ng/ul	100
19) Chrysene	21.19	228	25564	0.395	ng/ul	100
21) Benzo(b)fluoranthene	22.67	252	24351	0.426	ng/ul	100
22) Benzo(k)fluoranthene	22.72	252	24771	0.407	ng/ul	100
23) Benzo(a)pyrene	23.22	252	20533	0.423	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.47	276	25835	0.401	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.49	278	20796	0.400	ng/ul	100
26) Benzo(g,h,i)perylene	26.13	276	23185	0.407	ng/ul	99

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

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