

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN073020\
 Data File : BN011318.D
 Acq On : 31 Jul 2020 03:04
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.409

Quant Time: Jul 31 03:55:04 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 31 02:58:38 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	1097	0.40	ng/ul	0.00
2) Naphthalene-d8	10.21	136	4093	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	2652	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	6155	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	5973	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	6349	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	4774	0.46	ng/ul	0.00
14) Fluoranthene-d10	18.89	212	12332	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.26	128	4989	0.410	ng/ul	98
5) 2-Methylnaphthalene	11.89	142	3762	0.461	ng/ul	98
7) Acenaphthylene	13.82	152	5611	0.443	ng/ul	98
8) Acenaphthene	14.16	153	4179	0.415	ng/ul	98
9) Fluorene	15.16	166	4940	0.417	ng/ul	100
11) Pentachlorophenol	16.52	266	667	0.442	ng/ul	98
12) Phenanthrene	16.89	178	8426	0.413	ng/ul	99
13) Anthracene	16.99	178	7529	0.423	ng/ul	99
15) Fluoranthene	18.93	202	10324	0.411	ng/ul	98
17) Pyrene	19.29	202	10769	0.397	ng/ul	99
18) Benzo(a)anthracene	21.05	228	9772	0.426	ng/ul	99
19) Chrysene	21.10	228	10905	0.415	ng/ul	99
21) Benzo(b)fluoranthene	22.56	252	10584	0.399	ng/ul	98
22) Benzo(k)fluoranthene	22.61	252	11848	0.399	ng/ul	98
23) Benzo(a)pyrene	23.10	252	9924	0.431	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.26	276	11723	0.484	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.27	278	9267	0.481	ng/ul	100
26) Benzo(g,h,i)perylene	25.88	276	9560	0.437	ng/ul	99

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

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