

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN081020\
 Data File : BN011400.D
 Acq On : 10 Aug 2020 12:47
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.421

Quant Time: Aug 10 13:36:21 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 : Mon Aug 10 13:35:01 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	806	0.40	ng/ul	0.00
2) Naphthalene-d8	10.20	136	2308	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.09	164	1298	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.84	188	3090	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	2649	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	2608	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.80	152	2336	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.88	212	5058	0.34	ng/ul	0.00

Target Compounds						Ovalue
3) Naphthalene	10.25	128	2508	0.365	ng/ul	95
5) 2-Methylnaphthalene	11.88	142	1783	0.387	ng/ul	99
7) Acenaphthylene	13.80	152	2320	0.374	ng/ul	97
8) Acenaphthene	14.15	153	1836	0.373	ng/ul	98
9) Fluorene	15.15	166	2283	0.394	ng/ul	100
11) Pentachlorophenol	16.51	266	263	0.347	ng/ul	97
12) Phenanthrene	16.89	178	3726	0.364	ng/ul	98
13) Anthracene	16.98	178	3146	0.352	ng/ul	96
15) Fluoranthene	18.92	202	4305	0.341	ng/ul	99
17) Pyrene	19.28	202	3955	0.329	ng/ul	99
18) Benzo(a)anthracene	21.05	228	3277	0.322	ng/ul	98
19) Chrysene	21.10	228	4214	0.361	ng/ul	98
21) Benzo(b)fluoranthene	22.56	252	3807	0.350	ng/ul	91
22) Benzo(k)fluoranthene	22.60	252	4379	0.359	ng/ul	95
23) Benzo(a)pyrene	23.09	252	3491	0.369	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.24	276	5196	0.522	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.26	278	4121	0.520	ng/ul	96
26) Benzo(g,h,i)perylene	25.87	276	4772	0.531	ng/ul	95

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

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