

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072920\
 Data File : BN011293.D
 Acq On : 29 Jul 2020 15:20
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.413

Quant Time: Jul 29 15:51:00 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN072020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 29 11:20:12 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	1611	0.40	ng/ul	0.00
2) Naphthalene-d8	10.22	136	6019	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.11	164	3566	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	7348	0.40	ng/ul	0.00
16) Chrysene-d12	21.09	240	5086	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	4758	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.82	152	6377	0.43	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	13558	0.43	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.27	128	7516	0.417	ng/ul	99
5) 2-Methylnaphthalene	11.90	142	5045	0.428	ng/ul	99
7) Acenaphthylene	13.82	152	7257	0.420	ng/ul	98
8) Acenaphthene	14.17	153	5631	0.415	ng/ul	99
9) Fluorene	15.17	166	6375	0.414	ng/ul	99
11) Pentachlorophenol	16.53	266	757	0.482	ng/ul	96
12) Phenanthrene	16.90	178	10110	0.417	ng/ul	99
13) Anthracene	17.00	178	8846	0.443	ng/ul	98
15) Fluoranthene	18.93	202	11401	0.421	ng/ul	98
17) Pyrene	19.30	202	11348	0.418	ng/ul	99
18) Benzo(a)anthracene	21.07	228	8208	0.456	ng/ul	98
19) Chrysene	21.12	228	9456	0.395	ng/ul	99
21) Benzo(b)fluoranthene	22.59	252	7928	0.386	ng/ul	98
22) Benzo(k)fluoranthene	22.63	252	8696	0.369	ng/ul	98
23) Benzo(a)pyrene	23.13	252	7413	0.429	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.30	276	8498	0.417	ng/ul#	91
25) Dibenzo(a,h)anthracene	25.31	278	6796	0.442	ng/ul	96
26) Benzo(g,h,i)perylene	25.92	276	7553	0.373	ng/ul	99

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

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