

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN013051.D
 Acq On : 10 Dec 2020 17:47
 Operator : CG/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 121 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.439

Quant Time: Dec 10 18:17:25 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 10 10:29:27 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.37	152	899	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	3709	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.04	164	2025	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.80	188	4178	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	3028	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	2823	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	2361	0.43	ng/ul	0.00
14) Fluoranthene-d10	18.85	212	4580	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.19	128	4492	0.419	ng/ul	93
5) 2-Methylnaphthalene	11.82	142	3040	0.431	ng/ul	99
7) Acenaphthylene	13.76	152	4469	0.466	ng/ul	95
8) Acenaphthene	14.10	153	3376	0.429	ng/ul	98
9) Fluorene	15.10	166	3711	0.411	ng/ul	99
11) Pentachlorophenol	16.47	266	480	0.397	ng/ul	93
12) Phenanthrene	16.84	178	5912	0.422	ng/ul	98
13) Anthracene	16.94	178	5405	0.453	ng/ul	98
15) Fluoranthene	18.87	202	6732	0.404	ng/ul	99
17) Pyrene	19.24	202	6828	0.477	ng/ul	97
18) Benzo(a)anthracene	21.01	228	5486	0.438	ng/ul	99
19) Chrysene	21.06	228	6138	0.443	ng/ul	100
21) Benzo(b)fluoranthene	22.51	252	5672	0.434	ng/ul	95
22) Benzo(k)fluoranthene	22.55	252	5964	0.413	ng/ul	98
23) Benzo(a)pyrene	23.04	252	5216	0.430	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.18	276	6086	0.382	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.19	278	4854	0.386	ng/ul	96
26) Benzo(g,h,i)perylene	25.81	276	5266	0.366	ng/ul	97

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

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