

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120220\
 Data File : BN012800.D
 Acq On : 02 Dec 2020 02:04
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.417

Quant Time: Dec 02 02:34:16 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 02 00:53:21 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	673	0.40	ng/ul	0.00
2) Naphthalene-d8	10.19	136	2659	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.08	164	1509	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.84	188	3153	0.40	ng/ul	0.00
16) Chrysene-d12	21.05	240	3033	0.40	ng/ul	0.00
20) Perylene-d12	23.16	264	2996	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.79	152	1630	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.87	212	3237	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.24	128	3317	0.414	ng/ul	99
5) 2-Methylnaphthalene	11.86	142	2291	0.448	ng/ul	99
7) Acenaphthylene	13.80	152	3316	0.451	ng/ul	99
8) Acenaphthene	14.14	153	2515	0.406	ng/ul	98
9) Fluorene	15.14	166	2833	0.416	ng/ul	97
11) Pentachlorophenol	16.50	266	460	0.459	ng/ul	94
12) Phenanthrene	16.88	178	4476	0.408	ng/ul	98
13) Anthracene	16.97	178	4018	0.421	ng/ul	99
15) Fluoranthene	18.91	202	5548	0.438	ng/ul	98
17) Pyrene	19.27	202	5703	0.420	ng/ul	98
18) Benzo(a)anthracene	21.03	228	5581	0.513	ng/ul	99
19) Chrysene	21.09	228	6670	0.491	ng/ul	100
21) Benzo(b)fluoranthene	22.54	252	6612	0.480	ng/ul	94
22) Benzo(k)fluoranthene	22.58	252	7295	0.459	ng/ul	97
23) Benzo(a)pyrene	23.07	252	6303	0.486	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	25.23	276	8131	0.479	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.24	278	6499	0.477	ng/ul	95
26) Benzo(g,h,i)perylene	25.86	276	6986	0.458	ng/ul	98

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

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