

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080320\
 Data File : BN011364.D
 Acq On : 04 Aug 2020 14:46
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.417

Quant Time: Aug 04 15:19:55 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 03 10:06:56 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	1178	0.40	ng/ul	0.00
2) Naphthalene-d8	10.21	136	4266	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.09	164	2727	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	6202	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	6216	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	6057	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.80	152	4089	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.89	212	10888	0.36	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.26	128	4703	0.371	ng/ul	98
5) 2-Methylnaphthalene	11.88	142	3246	0.381	ng/ul	98
7) Acenaphthylene	13.81	152	4980	0.382	ng/ul	99
8) Acenaphthene	14.15	153	3886	0.375	ng/ul	98
9) Fluorene	15.15	166	4523	0.372	ng/ul	99
11) Pentachlorophenol	16.51	266	555	0.365	ng/ul	99
12) Phenanthrene	16.89	178	7652	0.372	ng/ul	99
13) Anthracene	16.98	178	6671	0.372	ng/ul	99
15) Fluoranthene	18.92	202	9264	0.366	ng/ul	99
17) Pyrene	19.28	202	9238	0.327	ng/ul	99
18) Benzo(a)anthracene	21.05	228	8308	0.348	ng/ul	99
19) Chrysene	21.10	228	9780	0.357	ng/ul	99
21) Benzo(b)fluoranthene	22.56	252	9365	0.370	ng/ul	99
22) Benzo(k)fluoranthene	22.60	252	10405	0.367	ng/ul	97
23) Benzo(a)pyrene	23.09	252	8234	0.375	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.25	276	9489	0.411	ng/ul	98
25) Dibenzo(a,h)anthracene	25.27	278	7594	0.413	ng/ul	99
26) Benzo(g,h,i)perylene	25.88	276	8151	0.390	ng/ul	99

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

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