

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120720\
 Data File : BN012958.D
 Acq On : 08 Dec 2020 03:08
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.430

Quant Time: Dec 08 03:39:39 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 : Mon Dec 07 12:24:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	882	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	3684	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	2075	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	4236	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	3345	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	3183	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	2212	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.86	212	4514	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.20	128	4218	0.396	ng/ul	94
5) 2-Methylnaphthalene	11.83	142	2834	0.405	ng/ul	98
7) Acenaphthylene	13.77	152	4057	0.413	ng/ul	96
8) Acenaphthene	14.12	153	3206	0.398	ng/ul	99
9) Fluorene	15.12	166	3603	0.390	ng/ul	98
11) Pentachlorophenol	16.47	266	499	0.407	ng/ul	97
12) Phenanthrene	16.85	178	5702	0.401	ng/ul	99
13) Anthracene	16.94	178	5058	0.418	ng/ul	97
15) Fluoranthene	18.88	202	6524	0.387	ng/ul	98
17) Pyrene	19.25	202	6653	0.421	ng/ul	100
18) Benzo(a)anthracene	21.02	228	5574	0.403	ng/ul	99
19) Chrysene	21.07	228	6268	0.410	ng/ul	99
21) Benzo(b)fluoranthene	22.51	252	6159	0.418	ng/ul	96
22) Benzo(k)fluoranthene	22.55	252	6594	0.405	ng/ul	95
23) Benzo(a)pyrene	23.04	252	5648	0.413	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.19	276	6516	0.363	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.20	278	5253	0.370	ng/ul	95
26) Benzo(g,h,i)perylene	25.82	276	5621	0.346	ng/ul	98

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

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