

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120620\
 Data File : BN012901.D
 Acq On : 06 Dec 2020 04:41
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.415

Quant Time: Dec 07 02:14:30 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 : Fri Dec 04 10:02:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	815	0.40	ng/ul	0.00
2) Naphthalene-d8	10.16	136	3115	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.06	164	1801	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.82	188	3834	0.40	ng/ul	0.00
16) Chrysene-d12	21.04	240	3300	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	3526	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	1880	0.41	ng/ul	-0.01
14) Fluoranthene-d10	18.86	212	4192	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.21	128	3532	0.392	ng/ul	96
5) 2-Methylnaphthalene	11.84	142	2354	0.398	ng/ul	99
7) Acenaphthylene	13.77	152	3376	0.396	ng/ul	97
8) Acenaphthene	14.12	153	2760	0.395	ng/ul	99
9) Fluorene	15.12	166	3111	0.388	ng/ul	98
11) Pentachlorophenol	16.48	266	446	0.402	ng/ul	96
12) Phenanthrene	16.86	178	5075	0.395	ng/ul	99
13) Anthracene	16.95	178	4472	0.409	ng/ul	97
15) Fluoranthene	18.89	202	6042	0.396	ng/ul	97
17) Pyrene	19.26	202	6155	0.395	ng/ul	99
18) Benzo(a)anthracene	21.02	228	5175	0.379	ng/ul	99
19) Chrysene	21.07	228	6191	0.410	ng/ul	100
21) Benzo(b)fluoranthene	22.52	252	6400	0.392	ng/ul	95
22) Benzo(k)fluoranthene	22.56	252	7328	0.406	ng/ul	94
23) Benzo(a)pyrene	23.05	252	6183	0.408	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.20	276	8277	0.416	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.21	278	6600	0.420	ng/ul	92
26) Benzo(g,h,i)perylene	25.83	276	7281	0.405	ng/ul	97

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

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