

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112818\
 Data File : BN003649.D
 Acq On : 28 Nov 2018 09:25
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.475

Quant Time: Nov 28 12:05:19 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 28 01:42:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	4113	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	15775	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	8207	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	19416	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	17860	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	15956	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	9422	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	22622	0.40	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.69	128	18064	0.408	ng/ul	100
5) 2-Methylnaphthalene	12.30	142	12579	0.407	ng/ul	99
7) Acenaphthylene	14.20	152	15376	0.416	ng/ul	100
8) Acenaphthene	14.54	153	13491	0.417	ng/ul	99
9) Fluorene	15.52	166	15467	0.410	ng/ul	99
11) Pentachlorophenol	16.87	266	1284	0.367	ng/ul	98
12) Phenanthrene	17.26	178	26733	0.420	ng/ul	99
13) Anthracene	17.36	178	21456	0.410	ng/ul	99
15) Fluoranthene	19.28	202	31411	0.417	ng/ul	98
17) Pyrene	19.64	202	30237	0.472	ng/ul	98
18) Benzo(a)anthracene	21.38	228	23135	0.410	ng/ul	99
19) Chrysene	21.44	228	26968	0.418	ng/ul	100
21) Benzo(b)fluoranthene	23.01	252	27669	0.419	ng/ul	98
22) Benzo(k)fluoranthene	23.06	252	26189	0.407	ng/ul	98
23) Benzo(a)pyrene	23.61	252	23397	0.409	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.07	276	24848	0.379	ng/ul	97
25) Dibenzo(a,h)anthracene	26.08	278	18722	0.355	ng/ul	98
26) Benzo(g,h,i)perylene	26.79	276	22132	0.389	ng/ul	97

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

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