

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120620\
 Data File : BN012916.D
 Acq On : 06 Dec 2020 14:17
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.416

Quant Time: Dec 07 03:04:06 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 02:16:16 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	817	0.40	ng/ul	0.00
2) Naphthalene-d8	10.16	136	3200	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.06	164	1880	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	4053	0.40	ng/ul	0.00
16) Chrysene-d12	21.04	240	3350	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	3457	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	1947	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.86	212	4359	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.21	128	3664	0.396	ng/ul	94
5) 2-Methylnaphthalene	11.84	142	2454	0.403	ng/ul	99
7) Acenaphthylene	13.77	152	3628	0.408	ng/ul	96
8) Acenaphthene	14.12	153	2935	0.402	ng/ul	99
9) Fluorene	15.12	166	3370	0.402	ng/ul	97
11) Pentachlorophenol	16.48	266	482	0.411	ng/ul	94
12) Phenanthrene	16.86	178	5395	0.397	ng/ul	99
13) Anthracene	16.95	178	4799	0.415	ng/ul	97
15) Fluoranthene	18.89	202	6316	0.391	ng/ul	97
17) Pyrene	19.26	202	6574	0.415	ng/ul	98
18) Benzo(a)anthracene	21.02	228	5421	0.391	ng/ul	98
19) Chrysene	21.07	228	6398	0.418	ng/ul	100
21) Benzo(b)fluoranthene	22.52	252	6414	0.401	ng/ul	95
22) Benzo(k)fluoranthene	22.56	252	7108	0.402	ng/ul	94
23) Benzo(a)pyrene	23.05	252	6141	0.414	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.20	276	7650	0.392	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.21	278	6136	0.398	ng/ul	92
26) Benzo(g,h,i)perylene	25.82	276	6701	0.380	ng/ul	97

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

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