

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120220\
 Data File : BN012774.D
 Acq On : 01 Dec 2020 10:11
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.415

Quant Time: Dec 01 17:06:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 01 17:04:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	607	0.40	ng/ul	0.00
2) Naphthalene-d8	10.19	136	2242	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.08	164	1346	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.84	188	2866	0.40	ng/ul	0.00
16) Chrysene-d12	21.05	240	2820	0.40	ng/ul	0.00
20) Perylene-d12	23.17	264	2771	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.79	152	1395	0.43	ng/ul	0.00
14) Fluoranthene-d10	18.88	212	2938	0.39	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.24	128	2877	0.426	ng/ul	99
5) 2-Methylnaphthalene	11.87	142	1953	0.453	ng/ul	100
7) Acenaphthylene	13.80	152	2843	0.433	ng/ul	99
8) Acenaphthene	14.14	153	2261	0.409	ng/ul	96
9) Fluorene	15.14	166	2521	0.415	ng/ul	99
11) Pentachlorophenol	16.50	266	400	0.439	ng/ul	97
12) Phenanthrene	16.88	178	4029	0.404	ng/ul	99
13) Anthracene	16.97	178	3600	0.415	ng/ul	98
15) Fluoranthene	18.91	202	5104	0.443	ng/ul	97
17) Pyrene	19.27	202	5223	0.414	ng/ul	100
18) Benzo(a)anthracene	21.04	228	4982	0.493	ng/ul	99
19) Chrysene	21.09	228	6270	0.497	ng/ul	98
21) Benzo(b)fluoranthene	22.54	252	5965	0.468	ng/ul	94
22) Benzo(k)fluoranthene	22.59	252	6836	0.465	ng/ul	95
23) Benzo(a)pyrene	23.08	252	5578	0.465	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.24	276	7522	0.479	ng/ul	98
25) Dibenzo(a,h)anthracene	25.25	278	5987	0.475	ng/ul	95
26) Benzo(g,h,i)perylene	25.87	276	6560	0.465	ng/ul	97

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

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