

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102120\
 Data File : BN012380.D
 Acq On : 21 Oct 2020 11:31
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.429

Quant Time: Oct 21 12:13:38 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 21 12:12:40 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	914	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	4042	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	2367	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	4633	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	3705	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	3851	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	2428	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	4536	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.41	128	4727	0.405	ng/ul	99
5) 2-Methylnaphthalene	12.04	142	3036	0.410	ng/ul	99
7) Acenaphthylene	13.95	152	3992	0.389	ng/ul	98
8) Acenaphthene	14.30	153	3452	0.398	ng/ul	98
9) Fluorene	15.29	166	3810	0.394	ng/ul	98
11) Pentachlorophenol	16.64	266	496	0.370	ng/ul	97
12) Phenanthrene	17.03	178	5779	0.391	ng/ul	99
13) Anthracene	17.12	178	4982	0.393	ng/ul	98
15) Fluoranthene	19.05	202	6268	0.363	ng/ul	98
17) Pyrene	19.41	202	6502	0.389	ng/ul	100
18) Benzo(a)anthracene	21.17	228	5142	0.375	ng/ul	97
19) Chrysene	21.22	228	6520	0.390	ng/ul	99
21) Benzo(b)fluoranthene	22.71	252	6573	0.423	ng/ul	96
22) Benzo(k)fluoranthene	22.76	252	6864	0.399	ng/ul	99
23) Benzo(a)pyrene	23.27	252	6080	0.402	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.52	276	7735	0.398	ng/ul#	55
25) Dibenzo(a,h)anthracene	25.54	278	6259	0.402	ng/ul	99
26) Benzo(g,h,i)perylene	26.18	276	7002	0.393	ng/ul	99

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

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