

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120319\
 Data File : BN008834.D
 Acq On : 03 Dec 2019 16:51
 Operator : JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.425

Quant Time: Dec 03 17:22:12 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 03 16:28:59 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	4438	0.40	ng/ul	0.00
2) Naphthalene-d8	10.34	136	18311	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	10222	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	20199	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	14721	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	12554	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	11547	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	24016	0.39	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.39	128	21291	0.386	ng/ul	99
5) 2-Methylnaphthalene	12.01	142	14738	0.382	ng/ul	99
7) Acenaphthylene	13.93	152	21474	0.380	ng/ul	100
8) Acenaphthene	14.27	153	15211	0.363	ng/ul	99
9) Fluorene	15.27	166	16994	0.345	ng/ul	99
11) Pentachlorophenol	16.62	266	2321	0.304	ng/ul	99
12) Phenanthrene	17.00	178	27357	0.402	ng/ul	99
13) Anthracene	17.09	178	26710	0.415	ng/ul	99
15) Fluoranthene	19.02	202	31632	0.378	ng/ul	99
17) Pyrene	19.39	202	32160	0.545	ng/ul	100
18) Benzo(a)anthracene	21.15	228	27222	0.460	ng/ul	98
19) Chrysene	21.20	228	25212	0.434	ng/ul	98
21) Benzo(b)fluoranthene	22.69	252	21451	0.414	ng/ul	86
22) Benzo(k)fluoranthene	22.73	252	21415	0.417	ng/ul#	84
23) Benzo(a)pyrene	23.23	252	19829	0.408	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	25.46	276	20744	0.365	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.47	278	16662	0.363	ng/ul#	87
26) Benzo(g,h,i)perylene	26.10	276	17420	0.366	ng/ul	93

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

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