

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120219\
 Data File : BN008809.D
 Acq On : 02 Dec 2019 11:41
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.422

Quant Time: Dec 02 13:11:00 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 : Mon Dec 02 12:46:36 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	3354	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	13632	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	8832	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	19426	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	16991	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	16392	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.94	152	9039	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	22899	0.39	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.39	128	16029	0.390	ng/ul	100
5) 2-Methylnaphthalene	12.01	142	11648	0.406	ng/ul	100
7) Acenaphthylene	13.93	152	17541	0.359	ng/ul	100
8) Acenaphthene	14.27	153	13149	0.363	ng/ul	100
9) Fluorene	15.26	166	15641	0.368	ng/ul	100
11) Pentachlorophenol	16.62	266	2087	0.284	ng/ul	100
12) Phenanthrene	17.00	178	26545	0.405	ng/ul	100
13) Anthracene	17.08	178	24560	0.397	ng/ul	100
15) Fluoranthene	19.02	202	30730	0.382	ng/ul	100
17) Pyrene	19.38	202	30772	0.451	ng/ul	100
18) Benzo(a)anthracene	21.14	228	29965	0.439	ng/ul	100
19) Chrysene	21.19	228	29192	0.435	ng/ul	100
21) Benzo(b)fluoranthene	22.67	252	28164	0.417	ng/ul	100
22) Benzo(k)fluoranthene	22.71	252	27695	0.413	ng/ul	100
23) Benzo(a)pyrene	23.22	252	25680	0.405	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.44	276	26469	0.357	ng/ul	100
25) Dibenzo(a,h)anthracene	25.45	278	21104	0.352	ng/ul	100
26) Benzo(g,h,i)perylene	26.08	276	22277	0.359	ng/ul	100

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

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