

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120219\
 Data File : BN008810.D
 Acq On : 02 Dec 2019 12:31
 Operator : JU
 Sample : K5851-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AA1

Quant Time: Dec 02 13:13:35 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	4639	0.40	ng/ul	0.00
2) Naphthalene-d8	10.34	136	18893	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	11216	0.40	ng/ul	0.00
10) Phenanthrene-d10	0.00	188	0	0.00	ng/ul	-16.95
16) Chrysene-d12	21.16	240	20479	0.40	ng/ul	0.00
20) Perylene-d12	23.32	264	20680	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	8268	0.27	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	20486	0.00	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.39	128	15477	0.272	ng/ul	99
5) 2-Methylnaphthalene	12.01	142	9033	0.227	ng/ul	100
7) Acenaphthylene	13.92	152	31808	0.513	ng/ul	99
8) Acenaphthene	14.27	153	2483	0.054	ng/ul	97
9) Fluorene	15.26	166	6764	0.125	ng/ul#	97
17) Pyrene	19.39	202	189123	2.302	ng/ul	95
18) Benzo(a)anthracene	21.14	228	139800	1.698	ng/ul	95
19) Chrysene	21.19	228	132469	1.638	ng/ul	98
21) Benzo(b)fluoranthene	22.68	252	238871	2.801	ng/ul	99
22) Benzo(k)fluoranthene	22.68	252	238871	2.824	ng/ul	98
23) Benzo(a)pyrene	23.23	252	123814	1.547	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.44	276	76158	0.814	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.45	278	17768	0.235	ng/ul#	76
26) Benzo(g,h,i)perylene	26.08	276	75550	0.965	ng/ul	100

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

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