

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN102419\
 Data File : BN008379.D
 Acq On : 24 Oct 2019 10:04
 Operator : JU
 Sample : K5235-04DL 5X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPB2DL

Quant Time: Oct 24 14:05:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 24 13:52:48 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	1341	0.40	ng/ul	0.00
2) Naphthalene-d8	10.34	136	6284	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	4075	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	8633	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	7151	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	4995	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	414	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	1258	0.04	ng/ul	0.00

Target Compounds				Qvalue		
7) Acenaphthylene	13.93	152	1642	0.085	ng/ul#	91
8) Acenaphthene	14.27	153	1723	0.114	ng/ul	96
9) Fluorene	15.27	166	3830	0.223	ng/ul	98
12) Phenanthrene	17.00	178	13296	0.526	ng/ul	100
13) Anthracene	17.09	178	13139	0.591	ng/ul	100
15) Fluoranthene	19.03	202	23960	0.718	ng/ul	98
17) Pyrene	19.39	202	26545	0.932	ng/ul	99
18) Benzo(a)anthracene	21.15	228	19344	0.744	ng/ul	99
19) Chrysene	21.20	228	23139	0.725	ng/ul	100
21) Benzo(b)fluoranthene	22.69	252	18599	1.144	ng/ul	96
22) Benzo(k)fluoranthene	22.73	252	17487	0.949	ng/ul	96
23) Benzo(a)pyrene	23.24	252	15346	0.897	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.48	276	16488	0.818	ng/ul#	91
25) Dibenzo(a,h)anthracene	25.49	278	12534	0.787	ng/ul	99
26) Benzo(g,h,i)perylene	26.13	276	14359	0.777	ng/ul	97

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

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