

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM031919\
 Data File : BM019236.D
 Acq On : 19 Mar 2019 14:40
 Operator : JU/SJ
 Sample : K1970-01DL 5X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P-1-N.WALLDL

Quant Time: Mar 19 15:42:13 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM031119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 19 12:12:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	217182	20.00	ng	0.00
21) Naphthalene-d8	10.43	136	979997	20.00	ng	0.00
39) Acenaphthene-d10	14.30	164	705046	20.00	ng	0.00
64) Phenanthrene-d10	17.06	188	1389344	20.00	ng	0.01
76) Chrysene-d12	21.25	240	1184242	20.00	ng	0.00
87) Perylene-d12	23.48	264	1108116	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.24	112	201337	15.01	ng	0.00
7) Phenol-d6	6.82	99	338089	17.24	ng	0.00
23) Nitrobenzene-d5	8.81	82	234573	11.31	ng	0.00
42) 2,4,6-Tribromophenol	15.80	330	202593	19.46	ng	0.00
45) 2-Fluorobiphenyl	12.91	172	684558	12.63	ng	0.00
79) Terphenyl-d14	19.69	244	985212	16.52	ng	0.00
Target Compounds						
22) Acetophenone	8.42	105	739439	27.303	ng	Qvalue # 61
31) Naphthalene	10.47	128	793660	15.365	ng	# 87
37) 2-Methylnaphthalene	12.10	142	3979698	106.026	ng	100
38) 1-Methylnaphthalene	12.33	142	3827985	103.467	ng	100
46) 1,1'-Biphenyl	13.13	154	1230729	19.988	ng	100
52) Acenaphthene	14.37	154	330818	7.409	ng	92
58) Fluorene	15.36	166	771192	12.992	ng	# 86
71) Phenanthrene	17.10	178	2645080	32.481	ng	# 96
78) Pyrene	19.49	202	555526	7.414	ng	# 86

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

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