

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM031919\
 Data File : BM019230.D
 Acq On : 19 Mar 2019 11:02
 Operator : JU/SJ
 Sample : K1970-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

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

Quant Time: Mar 19 12:27:57 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	7.65	152	242482	20.00	ng	0.00
21) Naphthalene-d8	10.45	136	1050161	20.00	ng	0.02
39) Acenaphthene-d10	14.33	164	897394	20.00	ng	0.04
64) Phenanthrene-d10	17.09	188	1161224	20.00	ng	0.04
76) Chrysene-d12	21.26	240	850022	20.00	ng	0.00
87) Perylene-d12	23.48	264	880461	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.24	112	1161869	77.60	ng	0.00
7) Phenol-d6	6.82	99	1884195	86.03	ng	0.00
23) Nitrobenzene-d5	8.82	82	1479161	66.58	ng	0.01
42) 2,4,6-Tribromophenol	15.84	330	1012401	76.41	ng	0.04
45) 2-Fluorobiphenyl	12.93	172	3512373	50.91	ng	0.02
79) Terphenyl-d14	19.70	244	3730529	87.13	ng	0.01
Target Compounds						
22) Acetophenone	8.43	105	3890287	134.046	ng	# 61
25) Isophorone	9.35	82	919830	21.698	ng	# 1
31) Naphthalene	10.50	128	3801034	68.671	ng	# 87
37) 2-Methylnaphthalene	12.14	142	18837610	468.337	ng	99
38) 1-Methylnaphthalene	12.36	142	17452885	440.220	ng	# 100
46) 1,1'-Biphenyl	13.16	154	6068004	77.427	ng	99
52) Acenaphthene	14.40	154	1534594	27.001	ng	# 85
58) Fluorene	15.39	166	3736728	49.459	ng	# 84
71) Phenanthrene	17.13	178	10334149	151.831	ng	# 95
78) Pyrene	19.50	202	2084134	38.752	ng	# 89

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

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