

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031519\
 Data File : BM019188.D
 Acq On : 15 Mar 2019 16:13
 Operator : JU/SJ
 Sample : K1899-03DL 5X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-4ADL

Manual Integrations
 APPROVED

Sohil
 3/18/2019 3:04:31 PM

Quant Time: Mar 16 02:27:27 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 11 17:30:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	201926	20.00	ng	-0.01
21) Naphthalene-d8	10.44	136	996205	20.00	ng	-0.01
39) Acenaphthene-d10	14.31	164	628616	20.00	ng	-0.01
64) Phenanthrene-d10	17.06	188	1396464	20.00	ng	-0.01
76) Chrysene-d12	21.26	240	1386297	20.00	ng	-0.01
87) Perylene-d12	23.50	264	1208571	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	158700	12.73	ng	0.00
7) Phenol-d6	6.84	99	147092	8.07	ng	0.00
23) Nitrobenzene-d5	8.82	82	375225	17.80	ng	-0.01
42) 2,4,6-Tribromophenol	15.81	330	260934	28.11	ng	-0.01
45) 2-Fluorobiphenyl	12.92	172	853011	17.65	ng	-0.01
79) Terphenyl-d14	19.70	244	1328145	19.02	ng	-0.01
Target Compounds						
17) 2-Methylphenol	8.10	107	36563	2.855	ng	92
27) 2,4-Dimethylphenol	9.63	122	72781m	5.389	ng	
31) Naphthalene	10.50	128	19006557	361.979	ng	# 89
37) 2-Methylnaphthalene	12.12	142	5838257	153.011	ng	100
38) 1-Methylnaphthalene	12.34	142	6064336	161.247	ng	100
46) 1,1'-Biphenyl	13.13	154	1059886	19.307	ng	99
52) Acenaphthene	14.37	154	1685196	42.328	ng	100
55) Dibenzofuran	14.71	168	1568423	24.428	ng	99
58) Fluorene	15.36	166	946166	17.878	ng	100
71) Phenanthrene	17.10	178	777746	9.502	ng	99
73) Carbazole	17.48	167	340411	4.501	ng	99

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

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