

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031519\
 Data File : BM019187.D
 Acq On : 15 Mar 2019 15:37
 Operator : JU/SJ
 Sample : K1899-02DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-2ADL

Manual Integrations
 APPROVED

Sohil
 3/18/2019 3:04:29 PM

Quant Time: Mar 16 02:25:36 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	197127	20.00	ng	-0.01
21) Naphthalene-d8	10.44	136	976149	20.00	ng	-0.01
39) Acenaphthene-d10	14.31	164	619161	20.00	ng	-0.01
64) Phenanthrene-d10	17.06	188	1370426	20.00	ng	-0.01
76) Chrysene-d12	21.26	240	1362126	20.00	ng	-0.01
87) Perylene-d12	23.50	264	1205056	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	144358	11.86	ng	0.00
7) Phenol-d6	6.83	99	136117	7.65	ng	-0.01
23) Nitrobenzene-d5	8.82	82	322482	15.62	ng	-0.01
42) 2,4,6-Tribromophenol	15.81	330	218900	23.95	ng	-0.01
45) 2-Fluorobiphenyl	12.92	172	747918	15.71	ng	-0.01
79) Terphenyl-d14	19.70	244	1161400	16.93	ng	-0.01
Target Compounds						
17) 2-Methylphenol	8.10	107	30611	2.448	ng	91
27) 2,4-Dimethylphenol	9.63	122	60947m	4.605	ng	
31) Naphthalene	10.50	128	17380777	337.817	ng	# 93
37) 2-Methylnaphthalene	12.12	142	4916200	131.493	ng	99
38) 1-Methylnaphthalene	12.34	142	5044532	136.887	ng	99
46) 1,1'-Biphenyl	13.13	154	886509	16.395	ng	99
52) Acenaphthene	14.37	154	1401431	35.738	ng	99
55) Dibenzofuran	14.71	168	1291106	20.416	ng	99
58) Fluorene	15.36	166	792875	15.210	ng	98
71) Phenanthrene	17.10	178	646108	8.044	ng	100
73) Carbazole	17.48	167	284933	3.839	ng	100

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

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