

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080718\
 Data File : BM016233.D
 Acq On : 08 Aug 2018 01:07
 Operator : SJ/JU
 Sample : J4297-03DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-3DL

Quant Time: Aug 08 03:18:43 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 07 16:15:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	56288	20.00	ng	0.00
21) Naphthalene-d8	10.96	136	298109	20.00	ng	0.00
38) Acenaphthene-d10	14.79	164	220418	20.00	ng	0.01
63) Phenanthrene-d10	17.53	188	447978	20.00	ng	0.02
75) Chrysene-d12	21.64	240	332824	20.00	ng	0.00
86) Perylene-d12	23.97	264	273828	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.67	112	56421	16.65	ng	0.00
7) Phenol-d6	7.30	99	74934	16.93	ng	0.00
23) Nitrobenzene-d5	9.33	82	124291	19.39	ng	0.00
41) 2,4,6-Tribromophenol	16.28	330	41273	14.76	ng	0.01
44) 2-Fluorobiphenyl	13.40	172	160452	8.86	ng	0.00
78) Terphenyl-d14	20.09	244	224101	14.09	ng	0.00
Target Compounds						
31) Naphthalene	11.02	128	1383363	91.689	ng	96
37) 2-Methylnaphthalene	12.64	142	6159722	609.460	ng	# 95
45) 1,1'-Biphenyl	13.62	154	812236	40.832	ng	96
70) Phenanthrene	17.57	178	995301	39.684	ng	# 94

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

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