

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080718\
 Data File : BM016236.D
 Acq On : 08 Aug 2018 02:55
 Operator : SJ/JU
 Sample : J4277-03 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OR-02-080318-B

Quant Time: Aug 08 04:25:20 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	8.13	152	48973	20.00	ng	0.00
21) Naphthalene-d8	10.95	136	212757	20.00	ng	0.00
38) Acenaphthene-d10	14.77	164	117099	20.00	ng	-0.01
63) Phenanthrene-d10	17.50	188	272805	20.00	ng	0.00
75) Chrysene-d12	21.64	240	315812	20.00	ng	0.00
86) Perylene-d12	23.98	264	271314	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.67	112	46115	15.64	ng	0.00
7) Phenol-d6	7.30	99	59438	15.44	ng	0.00
23) Nitrobenzene-d5	9.33	82	41297	9.03	ng	0.00
41) 2,4,6-Tribromophenol	16.26	330	19212	12.94	ng	-0.01
44) 2-Fluorobiphenyl	13.39	172	85888	8.92	ng	0.00
78) Terphenyl-d14	20.09	244	148011	9.81	ng	0.00
Target Compounds						
70) Phenanthrene	17.55	178	57448	3.761	ng	99
74) Fluoranthene	19.54	202	123286	7.236	ng	98
77) Pyrene	19.90	202	121940	6.443	ng	98
80) Benzo(a)anthracene	21.62	228	70865	3.643	ng	97
82) Chrysene	21.67	228	65349	3.502	ng	94
87) Benzo(b)fluoranthene	23.27	252	80275	5.025	ng	97
89) Benzo(a)pyrene	23.87	252	57233	3.727	ng	# 91
91) Benzo(g,h,i)perylene	27.10	276	39625	2.633	ng	96

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

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