

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM101818\
 Data File : BM017206.D
 Acq On : 18 Oct 2018 23:09
 Operator : MJ/SJ
 Sample : J5573-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 0692-KM-CA-AMD-COMP

Quant Time: Oct 19 13:56:18 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM101718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 17 14:23:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	290514	20.00	ng	0.00
21) Naphthalene-d8	10.45	136	1216714	20.00	ng	0.00
39) Acenaphthene-d10	14.31	164	709407	20.00	ng	0.00
64) Phenanthrene-d10	17.06	188	1554916	20.00	ng	-0.01
76) Chrysene-d12	21.25	240	1507812	20.00	ng	-0.01
87) Perylene-d12	23.46	264	1432509	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.28	112	839803	48.89	ng	0.00
7) Phenol-d6	6.85	99	673926	28.81	ng	-0.01
23) Nitrobenzene-d5	8.82	82	1668661	80.21	ng	-0.01
42) 2,4,6-Tribromophenol	15.80	330	1114415	116.16	ng	-0.01
45) 2-Fluorobiphenyl	12.93	172	4025653	75.50	ng	0.00
79) Terphenyl-d14	19.70	244	5306938	79.33	ng	0.00
Target Compounds						
10) Phenol	6.87	94	101249	4.021	ng	94
31) Naphthalene	10.49	128	160486	2.692	ng	99
50) Dimethylphthalate	13.77	163	227162	4.151	ng	99

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

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