

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062119\
 Data File : BM021079.D
 Acq On : 21 Jun 2019 12:33
 Operator : HP/JU
 Sample : K3309-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SS043MW010-190611

Quant Time: Jun 24 07:54:11 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM062019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 20 15:39:58 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	245970	20.00	ng	0.00
21) Naphthalene-d8	10.56	136	1046906	20.00	ng	0.00
39) Acenaphthene-d10	14.42	164	724185	20.00	ng	0.00
64) Phenanthrene-d10	17.16	188	1863551	20.00	ng	0.00
76) Chrysene-d12	21.34	240	1975441	20.00	ng	0.00
87) Perylene-d12	23.63	264	1986425	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	796726	58.65	ng	0.00
7) Phenol-d6	6.95	99	707691	35.77	ng	0.00
23) Nitrobenzene-d5	8.94	82	1798020	89.51	ng	0.00
42) 2,4,6-Tribromophenol	15.91	330	1935647	142.44	ng	0.00
45) 2-Fluorobiphenyl	13.03	172	5309743	90.11	ng	0.00
79) Terphenyl-d14	19.79	244	9909220	93.69	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.32	88	25711	5.024	ng	Qvalue # 92
13) 1,4-Dichlorobenzene	7.81	146	47370	2.300	ng	95
52) Acenaphthene	14.48	154	422170	9.134	ng	100
58) Fluorene	15.47	166	425232	6.902	ng	100

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

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