

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062219\
 Data File : BM021095.D
 Acq On : 22 Jun 2019 11:22
 Operator : HP/JU
 Sample : K3309-01DL 20X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SS043MW009-190610DL

Quant Time: Jun 24 06:32:02 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	7.77	152	240247	20.00	ng	0.00
21) Naphthalene-d8	10.56	136	1045104	20.00	ng	0.00
39) Acenaphthene-d10	14.42	164	719034	20.00	ng	0.00
64) Phenanthrene-d10	17.16	188	1746703	20.00	ng	0.00
76) Chrysene-d12	21.34	240	1460418	20.00	ng	0.00
87) Perylene-d12	23.62	264	1411858	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	33840	2.55	ng	0.00
7) Phenol-d6	6.95	99	30208	1.56	ng	0.00
23) Nitrobenzene-d5	8.94	82	76273	3.80	ng	0.00
42) 2,4,6-Tribromophenol	15.91	330	70106	5.20	ng	0.00
45) 2-Fluorobiphenyl	13.03	172	243424	4.16	ng	0.00
79) Terphenyl-d14	19.79	244	343408	4.39	ng	0.00
Target Compounds						
12) 1,3-Dichlorobenzene	7.66	146	54237	2.742	ng	97
13) 1,4-Dichlorobenzene	7.81	146	206219	10.253	ng	99
14) 1,2-Dichlorobenzene	8.13	146	1627017	82.542	ng	99

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

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