

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081719\
 Data File : BG042277.D
 Acq On : 17 Aug 2019 2:19
 Operator : HP/JU
 Sample : K4264-16
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 REACTOR-3-G1-G4-(3)

Quant Time: Aug 17 05:43:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG072719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 31 08:57:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	56197	20.00	ng	-0.03
21) Naphthalene-d8	10.83	136	212646	20.00	ng	-0.04
39) Acenaphthene-d10	14.67	164	156058	20.00	ng	-0.03
64) Phenanthrene-d10	17.43	188	388902	20.00	ng	-0.03
76) Chrysene-d12	21.70	240	448539	20.00	ng	-0.03
87) Perylene-d12	24.96	264	518049	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	387449	134.14	ng	-0.02
7) Phenol-d6	7.17	99	469333	120.53	ng	-0.02
23) Nitrobenzene-d5	9.18	82	322723	104.44	ng	-0.03
42) 2,4,6-Tribromophenol	16.16	330	340508	192.09	ng	-0.03
45) 2-Fluorobiphenyl	13.29	172	995367	112.49	ng	-0.04
79) Terphenyl-d14	20.04	244	1963302	100.16	ng	-0.03

Target Compounds Qvalue

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

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