

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051721\
 Data File : BF123961.D
 Acq On : 17 May 2021 13:59
 Operator : JU/CG
 Sample : M2383-07DL 2X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 KTW-01DL

Quant Time: May 17 14:30:02 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 14 19:05:57 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.904	152	41738	20.00	ng	0.00
21) Naphthalene-d8	8.186	136	166675	20.00	ng	# 0.00
39) Acenaphthene-d10	9.939	164	94817	20.00	ng	0.00
64) Phenanthrene-d10	11.427	188	182280	20.00	ng	# 0.00
76) Chrysene-d12	14.068	240	149379	20.00	ng	0.00
86) Perylene-d12	15.562	264	131586	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	58438	19.66	ng	0.00
7) Phenol-d6	6.510	99	42046	10.90	ng	-0.01
23) Nitrobenzene-d5	7.463	82	143513	37.10	ng	0.00
42) 2,4,6-Tribromophenol	10.727	330	68879	60.48	ng	0.00
45) 2-Fluorobiphenyl	9.263	172	275196	35.72	ng	0.00
79) Terphenyl-d14	13.015	244	314003	31.84	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.210	128	361848	36.68	ng	98
37) 2-Methylnaphthalene	8.898	142	39748	5.95	ng	100
38) 1-Methylnaphthalene	8.998	142	24775	3.97	ng	96
49) Acenaphthylene	9.804	152	30467	3.11	ng	# 96
55) Dibenzofuran	10.145	168	28960	3.01	ng	97
58) Fluorene	10.486	166	30536	4.21	ng	83
71) Phenanthrene	11.451	178	72480	6.27	ng	98

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

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