

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072021\
 Data File : BF124645.D
 Acq On : 21 Jul 2021 00:53
 Operator : JU/CG
 Sample : M3089-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PAR-N2-(A-D)-(0-12)

Quant Time: Jul 21 03:14:22 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF070721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 20 13:36:28 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.904	152	7338	20.000	ng	-0.01
21) Naphthalene-d8	8.186	136	28965	20.000	ng	#-0.02
39) Acenaphthene-d10	9.945	164	14999	20.000	ng	-0.01
64) Phenanthrene-d10	11.433	188	19816	20.000	ng	-0.01
76) Chrysene-d12	14.068	240	11064	20.000	ng	-0.02
86) Perylene-d12	15.557	264	10592	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	43180	101.952	ng	-0.01
7) Phenol-d6	6.522	99	50821	98.638	ng	-0.02
23) Nitrobenzene-d5	7.463	82	31579	68.824	ng	-0.02
42) 2,4,6-Tribromophenol	10.733	330	10106	86.583	ng	-0.02
45) 2-Fluorobiphenyl	9.263	172	67894	66.008	ng	-0.02
79) Terphenyl-d14	13.021	244	38519	58.461	ng	-0.01
Target Compounds						
						Qvalue
71) Phenanthrene	11.457	178	2662	2.462	ng	96
75) Fluoranthene	12.645	202	2390	2.191	ng	89
78) Pyrene	12.874	202	2059	2.240	ng	# 87

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

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