

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061821\
 Data File : BF124377.D
 Acq On : 18 Jun 2021 16:52
 Operator : JU/CG
 Sample : M2687-15
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 17-B6(116-120)

Manual Integrations
 APPROVED

mohammad
 6/21/2021 11:38:27 AM

Quant Time: Jun 18 17:21:06 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 15 20:09:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.786	152	28217	20.000	ng	0.00
21) Naphthalene-d8	8.069	136	115450	20.000	ng	# 0.00
39) Acenaphthene-d10	9.822	164	64407	20.000	ng	0.00
64) Phenanthrene-d10	11.310	188	136009	20.000	ng	0.00
76) Chrysene-d12	13.957	240	129268	20.000	ng	0.00
86) Perylene-d12	15.415	264	97410	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	238745	127.368	ng	0.00
7) Phenol-d6	6.434	99	270060	107.185	ng	-0.01
23) Nitrobenzene-d5	7.351	82	240664	82.677	ng	0.00
42) 2,4,6-Tribromophenol	10.616	330	128861	141.219	ng	0.00
45) 2-Fluorobiphenyl	9.145	172	459530	89.782	ng	0.00
79) Terphenyl-d14	12.904	244	676519	71.853	ng	0.00
Target Compounds						
77) Benzidine	12.651	184	21743m	10.667	ng	Qvalue

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

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