

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121621\
 Data File : BF126208.D
 Acq On : 16 Dec 2021 13:08
 Operator : JU/CG
 Sample : M4919-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FDR-MW-2-20211202

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 12/17/2021
 Supervised By :Jagrut Upadhyay 12/17/2021

Quant Time: Dec 16 13:54:10 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 16 12:22:48 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.816	152	179769	20.000	ng	0.00
21) Naphthalene-d8	8.098	136	754662	20.000	ng	# 0.00
39) Acenaphthene-d10	9.851	164	347049	20.000	ng	0.00
64) Phenanthrene-d10	11.333	188	457873m	20.000	ng	0.00
76) Chrysene-d12	13.968	240	335023	20.000	ng	# 0.00
86) Perylene-d12	15.415	264	341115	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	689628	56.494	ng	0.00
7) Phenol-d6	6.428	99	533735	34.434	ng	-0.02
23) Nitrobenzene-d5	7.375	82	1267903	90.992	ng	0.00
42) 2,4,6-Tribromophenol	10.639	330	338429	121.342	ng	0.00
45) 2-Fluorobiphenyl	9.175	172	1686569	85.234	ng	0.00
79) Terphenyl-d14	12.927	244	1108939	57.532	ng	0.00

Target Compounds Qvalue

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

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