

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111321\
 Data File : BF126172.D
 Acq On : 13 Nov 2021 17:13
 Operator : JU/CG
 Sample : M4593-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WASTE-108

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 11/15/2021
 Supervised By :mohammad ahmed 11/17/2021

Quant Time: Nov 15 09:58:59 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 15 08:09:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	104221	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	397798	20.000	ng	0.00
39) Acenaphthene-d10	9.869	164	200893	20.000	ng	0.00
64) Phenanthrene-d10	11.351	188	369643	20.000	ng	# 0.00
76) Chrysene-d12	13.992	240	288983	20.000	ng	#-0.01
86) Perylene-d12	15.457	264	331054	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.457	112	627536	103.401	ng	0.01
7) Phenol-d6	6.463	99	601573	70.316	ng	0.00
23) Nitrobenzene-d5	7.392	82	679285	80.532	ng	0.00
42) 2,4,6-Tribromophenol	10.657	330	389641	169.244	ng	0.00
45) 2-Fluorobiphenyl	9.192	172	1325634	89.172	ng	0.00
79) Terphenyl-d14	12.945	244	1687137	94.143	ng	0.00
Target Compounds						Qvalue
32) Benzoic acid	7.822	122	1641m	9.529	ng	

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

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