

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030422\
 Data File : BF127206.D
 Acq On : 04 Mar 2022 17:45
 Operator : CG\JU
 Sample : N1793-09
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-3

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/07/2022
 Supervised By : Jagrut Upadhyay 03/07/2022

Quant Time: Mar 05 02:21:22 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 05 02:16:58 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	93230	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	356377	20.000	ng	# 0.00
39) Acenaphthene-d10	9.916	164	191291	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	350296	20.000	ng	# 0.00
76) Chrysene-d12	14.057	240	257509	20.000	ng	# 0.00
86) Perylene-d12	15.545	264	209486	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	292728	48.529	ng	0.00
7) Phenol-d6	6.504	99	225195m	27.667	ng	-0.01
23) Nitrobenzene-d5	7.445	82	703992	89.165	ng	0.00
42) 2,4,6-Tribromophenol	10.716	330	333700	151.513	ng	0.00
45) 2-Fluorobiphenyl	9.239	172	1137599	87.548	ng	0.00
79) Terphenyl-d14	13.004	244	1316094	75.132	ng	0.00

Target Compounds Qvalue

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

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