

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF093021\
 Data File : BF125729.D
 Acq On : 30 Sep 2021 12:57
 Operator : JU/CG
 Sample : M3951-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-20210924

Manual Integrations
 APPROVED

mohammad
 10/1/2021 11:28:46 AM

Quant Time: Sep 30 13:42:10 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 24 15:38:32 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.886	152	216630	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	875823	20.000	ng	# 0.00
39) Acenaphthene-d10	9.922	164	458143	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	791566	20.000	ng	# 0.00
76) Chrysene-d12	14.039	240	579616	20.000	ng	0.00
86) Perylene-d12	15.515	264	444387	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	1584196	120.485	ng	0.01
7) Phenol-d6	6.510	99	1709043	96.681	ng	0.00
23) Nitrobenzene-d5	7.445	82	1307400	104.121	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	638273	149.423	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	2372304	94.543	ng	0.00
79) Terphenyl-d14	12.992	244	2594883	77.529	ng	0.00
Target Compounds						
32) Benzoic acid	7.851	122	3746m	8.963	ng	Qvalue
77) Benzidine	12.733	184	614478	40.355	ng	96

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF093021\
Data File : BF125729.D
Acq On : 30 Sep 2021 12:57
Operator : JU/CG
Sample : M3951-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-20210924

Manual Integrations
APPROVED

mohammad
10/1/2021 11:28:46 AM

Quant Time: Sep 30 13:42:10 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 24 15:38:32 2021
Response via : Initial Calibration

