

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102721\
 Data File : BF125964.D
 Acq On : 27 Oct 2021 15:30
 Operator : JU/CG
 Sample : M4325-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RED-BOX-2-1934

Manual Integrations
 APPROVED

mohammad
 10/29/2021 2:22:24 PM

Quant Time: Oct 28 08:58:24 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 25 18:49:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.857	152	170327	20.000	ng	-0.01
21) Naphthalene-d8	8.139	136	690999	20.000	ng	#-0.02
39) Acenaphthene-d10	9.892	164	391538	20.000	ng	-0.02
64) Phenanthrene-d10	11.374	188	755156	20.000	ng	-0.02
76) Chrysene-d12	14.010	240	518446	20.000	ng	-0.02
86) Perylene-d12	15.474	264	453271	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	1350658	118.116	ng	0.00
7) Phenol-d6	6.481	99	1541937	98.313	ng	-0.01
23) Nitrobenzene-d5	7.416	82	1241732	92.822	ng	-0.02
42) 2,4,6-Tribromophenol	10.680	330	519255	152.131	ng	-0.01
45) 2-Fluorobiphenyl	9.216	172	1991371	89.943	ng	-0.02
79) Terphenyl-d14	12.968	244	2541263	80.178	ng	-0.01
Target Compounds						
32) Benzoic acid	7.833	122	33592m	8.396	ng	Qvalue

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

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