

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102320\
 Data File : BF122110.D
 Acq On : 23 Oct 2020 20:53
 Operator : JU/CG
 Sample : L4487-03 10X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 D-89

Manual Integrations
 APPROVED

mohammad
 10/26/2020 10:35:29 AM

Quant Time: Oct 23 23:37:32 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 21 12:04:42 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.722	152	102479	20.00	ng	0.00
21) Naphthalene-d8	8.010	136	419243	20.00	ng	0.00
39) Acenaphthene-d10	9.757	164	205605	20.00	ng	0.00
64) Phenanthrene-d10	11.245	188	338260	20.00	ng	0.00
76) Chrysene-d12	13.892	240	239604	20.00	ng	0.00
86) Perylene-d12	15.333	264	241941	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	4953m	0.71	ng	0.07
7) Phenol-d6	6.451	99	7396	0.83	ng	0.07
23) Nitrobenzene-d5	7.328	82	5152	0.63	ng	0.02
42) 2,4,6-Tribromophenol	10.569	330	755	0.30	ng	0.01
45) 2-Fluorobiphenyl	9.086	172	12243	0.88	ng	0.00
79) Terphenyl-d14	12.827	244	10831	0.86	ng	0.00

Target Compounds	Qvalue

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

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