

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072820\
 Data File : BF121082.D
 Acq On : 28 Jul 2020 21:54
 Operator : JU/CG
 Sample : L3439-13
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20200443-3002C

Manual Integrations
 APPROVED

mohammad
 7/29/2020 11:47:39 AM

Quant Time: Jul 29 06:33:01 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 16 14:20:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	156163	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	596936	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	319131	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	635977	20.00	ng	0.00
76) Chrysene-d12	13.97	240	630796	20.00	ng	0.00
86) Perylene-d12	15.42	264	648632	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	1212204	127.25	ng	0.02
7) Phenol-d6	6.45	99	1370267	111.36	ng	0.00
23) Nitrobenzene-d5	7.38	82	1008138	97.72	ng	0.00
42) 2,4,6-Tribromophenol	10.64	330	532225	152.36	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	1807805	84.58	ng	0.00
79) Terphenyl-d14	12.92	244	2031324	70.00	ng	0.00
Target Compounds						
10) Phenol	6.46	94	75603	5.585	ng	96
32) Benzoic acid	7.83	122	22031m	3.423	ng	
50) Dimethylphthalate	9.56	163	168452	7.317	ng	99

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

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