

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101420\
 Data File : BF121961.D
 Acq On : 14 Oct 2020 16:59
 Operator : JU/CG
 Sample : L4380-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B-BRICK-2

Manual Integrations
 APPROVED

mohammad
 10/15/2020 10:47:13 AM

Quant Time: Oct 14 17:42:34 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 12 16:22:33 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.739	152	111789	20.00	ng	0.00
21) Naphthalene-d8	8.028	136	442952	20.00	ng	0.00
39) Acenaphthene-d10	9.775	164	225904	20.00	ng	-0.01
64) Phenanthrene-d10	11.263	188	414364	20.00	ng	-0.01
76) Chrysene-d12	13.904	240	292692	20.00	ng	-0.01
86) Perylene-d12	15.339	264	239278	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	600693	79.31	ng	0.00
7) Phenol-d6	6.387	99	762249	78.18	ng	-0.01
23) Nitrobenzene-d5	7.310	82	516581	59.81	ng	-0.01
42) 2,4,6-Tribromophenol	10.569	330	194168	69.48	ng	-0.01
45) 2-Fluorobiphenyl	9.098	172	911953	59.40	ng	-0.01
79) Terphenyl-d14	12.845	244	911783	59.37	ng	-0.02
Target Compounds						Qvalue
50) Dimethylphthalate	9.498	163	150629	9.12	ng	100
87) Indeno(1,2,3-cd)pyrene	16.745	276	130022	8.37	ng	98
88) Benzo(b)fluoranthene	14.927	252	43962	2.72	ng	99
89) Benzo(k)fluoranthene	14.951	252	40119m	2.61	ng	
90) Benzo(a)pyrene	15.280	252	43636	3.12	ng	# 94
91) Dibenzo(a,h)anthracene	16.751	278	94679	7.40	ng	99
92) Benzo(g,h,i)perylene	17.168	276	128929	10.07	ng	99

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

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