

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101320\
 Data File : BF121948.D
 Acq On : 13 Oct 2020 20:21
 Operator : JU/CG
 Sample : L4372-01 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-HADLEY

Manual Integrations
 APPROVED

mohammad
 10/14/2020 12:53:38 PM

Quant Time: Oct 14 02:07:55 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	6.745	152	93840	20.00	ng	0.00
21) Naphthalene-d8	8.028	136	330074	20.00	ng	0.00
39) Acenaphthene-d10	9.775	164	141538	20.00	ng	-0.01
64) Phenanthrene-d10	11.263	188	252876	20.00	ng	-0.01
76) Chrysene-d12	13.904	240	274938	20.00	ng	-0.01
86) Perylene-d12	15.339	264	262049	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	102073	16.05	ng	0.00
7) Phenol-d6	6.387	99	126001	15.39	ng	-0.01
23) Nitrobenzene-d5	7.310	82	80163	12.46	ng	-0.01
42) 2,4,6-Tribromophenol	10.569	330	23820	13.60	ng	-0.01
45) 2-Fluorobiphenyl	9.098	172	146721	15.25	ng	-0.01
79) Terphenyl-d14	12.839	244	145456	10.08	ng	-0.02
Target Compounds						Qvalue
75) Fluoranthene	12.474	202	48713	3.28	ng	97
78) Pyrene	12.704	202	44657	2.13	ng	97
88) Benzo(b)fluoranthene	14.927	252	45607m	2.57	ng	

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

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