

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF021120\
 Data File : BF118899.D
 Acq On : 11 Feb 2020 22:29
 Operator : CG/JU
 Sample : L1367-01 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-03-021020

Manual Integrations
 APPROVED

mohammad
 2/12/2020 3:51:13 PM

Quant Time: Feb 12 07:47:04 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF020320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 03 16:26:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	265099	20.00	ng	-0.01
21) Naphthalene-d8	8.09	136	1054729	20.00	ng	-0.02
39) Acenaphthene-d10	9.84	164	436887	20.00	ng	-0.02
64) Phenanthrene-d10	11.33	188	777310	20.00	ng	-0.01
76) Chrysene-d12	13.96	240	536875	20.00	ng	-0.02
87) Perylene-d12	15.42	264	611871	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	781344	55.79	ng	0.00
7) Phenol-d6	6.44	99	926898	53.32	ng	-0.01
23) Nitrobenzene-d5	7.36	82	603000	34.08	ng	-0.02
42) 2,4,6-Tribromophenol	10.63	330	245088	46.88	ng	-0.02
45) 2-Fluorobiphenyl	9.16	172	1316423	53.87	ng	-0.02
79) Terphenyl-d14	12.91	244	1012520	38.74	ng	-0.02
Target Compounds						
71) Phenanthrene	11.35	178	85643	2.239	ng	95
75) Fluoranthene	12.54	202	153000	3.664	ng	99
78) Pyrene	12.77	202	147423	4.003	ng	97
83) Chrysene	13.99	228	69173	2.157	ng	95
88) Benzo(b)fluoranthene	15.00	252	92775m	2.444	ng	
90) Benzo(a)pyrene	15.35	252	66167	2.036	ng	# 92

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

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