

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030221\
 Data File : BF123351.D
 Acq On : 2 Mar 2021 23:05
 Operator : JU/CG
 Sample : M1525-01 10X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-01-030121

Manual Integrations
 APPROVED

mohammad
 3/3/2021 11:01:51 AM

Quant Time: Mar 02 23:53:59 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF021921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 19 16:08:53 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.851	152	130926	20.00	ng	0.00
21) Naphthalene-d8	8.134	136	538802	20.00	ng	#-0.01
39) Acenaphthene-d10	9.892	164	276942	20.00	ng	-0.01
64) Phenanthrene-d10	11.386	188	479111	20.00	ng	0.00
76) Chrysene-d12	14.039	240	295820	20.00	ng	0.00
86) Perylene-d12	15.521	264	289766	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	63535	6.99	ng	0.00
7) Phenol-d6	6.487	99	81918	6.63	ng	0.00
23) Nitrobenzene-d5	7.410	82	54312	4.61	ng	-0.02
42) 2,4,6-Tribromophenol	10.686	330	17343	6.15	ng	-0.01
45) 2-Fluorobiphenyl	9.210	172	90212	4.68	ng	-0.01
79) Terphenyl-d14	12.974	244	81705	5.31	ng	-0.01
Target Compounds						Qvalue
71) Phenanthrene	11.410	178	171247	6.04	ng	99
75) Fluoranthene	12.604	202	138347	4.61	ng	99
78) Pyrene	12.833	202	151779	6.57	ng	99
81) Benzo(a)anthracene	14.027	228	41286	2.05	ng	# 90
83) Chrysene	14.063	228	52208m	2.64	ng	
90) Benzo(a)pyrene	15.457	252	36418	2.01	ng	# 79

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

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