

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF073021\
 Data File : BF124883.D
 Acq On : 30 Jul 2021 18:59
 Operator : JU/CG
 Sample : M3238-01 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 282

Manual Integrations
 APPROVED

mohammad
 8/2/2021 3:44:54 PM

Quant Time: Jul 31 00:07:26 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 22 15:28:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	7745	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	32427	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	18189	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	28646	20.000	ng	0.00
76) Chrysene-d12	14.039	240	14442	20.000	ng	0.00
86) Perylene-d12	15.521	264	12778	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	22883	50.025	ng	0.00
7) Phenol-d6	6.498	99	28430	49.577	ng	0.00
23) Nitrobenzene-d5	7.433	82	19984	36.679	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	8090	51.802	ng	0.00
45) 2-Fluorobiphenyl	9.227	172	43582	35.192	ng	0.00
79) Terphenyl-d14	12.986	244	33049	39.226	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	12.615	202	3400	2.163	ng	83
78) Pyrene	12.845	202	3333	2.915	ng	# 90
84) Bis(2-ethylhexyl)phtha...	14.009	149	12126	15.033	ng	99
88) Benzo(b)fluoranthene	15.080	252	2577m	2.864	ng	

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

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