

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101821\
 Data File : BF125889.D
 Acq On : 18 Oct 2021 21:29
 Operator : JU/CG
 Sample : M4259-04 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-1-BS

Manual Integrations
 APPROVED

mohammad
 10/19/2021 12:10:01 PM

Quant Time: Oct 19 09:09:04 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF100721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 18 12:43:03 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.822	152	177958	20.00	ng	0.00
21) Naphthalene-d8	8.104	136	700570	20.00	ng	# 0.00
39) Acenaphthene-d10	9.857	164	314468	20.00	ng	0.00
64) Phenanthrene-d10	11.339	188	486579	20.00	ng	0.00
76) Chrysene-d12	13.974	240	411420	20.00	ng	0.00
86) Perylene-d12	15.421	264	275102	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	433389	39.78	ng	0.00
7) Phenol-d6	6.445	99	512143	36.52	ng	-0.01
23) Nitrobenzene-d5	7.381	82	283789	25.18	ng	-0.01
42) 2,4,6-Tribromophenol	10.639	330	106501	34.30	ng	0.00
45) 2-Fluorobiphenyl	9.175	172	556293	30.51	ng	0.00
79) Terphenyl-d14	12.922	244	508673	18.83	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	12.545	202	101820	4.33	ng	94
78) Pyrene	12.774	202	103927	2.49	ng	98
88) Benzo(b)fluoranthene	15.004	252	59526m	3.83	ng	
90) Benzo(a)pyrene	15.357	252	41530	3.02	ng	# 94
92) Benzo(g,h,i)perylene	17.286	276	41442	2.34	ng	95

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

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