

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092321\
 Data File : BF125606.D
 Acq On : 23 Sep 2021 22:55
 Operator : JU/CG
 Sample : M3730-05
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WALL-BRICK-FLOOR

Quant Time: Sep 24 04:19:49 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 21 19:21:36 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.893	152	219100	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	903389	20.000	ng	# 0.00
39) Acenaphthene-d10	9.928	164	491141	20.000	ng	-0.01
64) Phenanthrene-d10	11.416	188	870823	20.000	ng	# 0.00
76) Chrysene-d12	14.057	240	718295	20.000	ng	0.00
86) Perylene-d12	15.533	264	626844	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	1545569	111.552	ng	0.01
7) Phenol-d6	6.516	99	1694430	94.852	ng	0.00
23) Nitrobenzene-d5	7.451	82	1227382	86.657	ng	-0.01
42) 2,4,6-Tribromophenol	10.722	330	645295	134.141	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	2453118	91.255	ng	-0.01
79) Terphenyl-d14	13.010	244	2736862	61.528	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.198	128	358989	7.738	ng	97
37) 2-Methylnaphthalene	8.887	142	85584	2.723	ng	98
52) Acenaphthene	9.963	154	108822	3.816	ng	98
55) Dibenzofuran	10.134	168	122623	2.898	ng	93
58) Fluorene	10.475	166	85005	2.612	ng	100
71) Phenanthrene	11.439	178	214673	4.389	ng	100

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

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