

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102623\
 Data File : BF135972.D
 Acq On : 26 Oct 2023 19:52
 Operator : CG\JU
 Sample : 04685-03DL 10X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PP-20230928-B4A-7.5-15DL

Quant Time: Oct 27 01:19:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 26 03:58:01 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.828	152	159064	20.000	ng	0.00
21) Naphthalene-d8	8.104	136	532973	20.000	ng	# 0.00
39) Acenaphthene-d10	9.869	164	272789	20.000	ng	0.00
64) Phenanthrene-d10	11.357	188	412430	20.000	ng	0.00
76) Chrysene-d12	14.004	240	347495	20.000	ng	0.00
86) Perylene-d12	15.492	264	274359	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	104265	10.122	ng	-0.01
7) Phenol-d6	6.440	99	132725	10.349	ng	-0.02
23) Nitrobenzene-d5	7.381	82	79757	18.339	ng	-0.01
42) 2,4,6-Tribromophenol	10.651	330	17665	7.398	ng	0.00
45) 2-Fluorobiphenyl	9.186	172	144273	8.469	ng	0.00
79) Terphenyl-d14	12.951	244	131518	6.323	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.128	128	384356	14.916	ng	# 96
37) 2-Methylnaphthalene	8.828	142	1078241	63.244	ng	99
38) 1-Methylnaphthalene	8.922	142	661494	41.523	ng	100
46) 1,1'-Biphenyl	9.286	154	128129	6.199	ng	98
52) Acenaphthene	9.898	154	42835	2.771	ng	# 82
58) Fluorene	10.416	166	106679	6.385	ng	# 89
71) Phenanthrene	11.380	178	264412	12.672	ng	96

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

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