

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072622\
 Data File : BF129363.D
 Acq On : 26 Jul 2022 19:24
 Operator : CG\JU
 Sample : N3886-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WH-1

Quant Time: Jul 27 00:40:53 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 25 17:24:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	218907	20.000	ng	0.00
21) Naphthalene-d8	8.163	136	842183	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	411987	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	577304	20.000	ng	# 0.00
76) Chrysene-d12	14.057	240	413084	20.000	ng	-0.01
86) Perylene-d12	15.545	264	456340	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	1688899	125.680	ng	0.00
7) Phenol-d6	6.510	99	2121737	125.614	ng	0.00
23) Nitrobenzene-d5	7.445	82	1332901	86.396	ng	0.00
42) 2,4,6-Tribromophenol	10.716	330	434891	109.795	ng	0.00
45) 2-Fluorobiphenyl	9.239	172	2115338	79.428	ng	0.00
79) Terphenyl-d14	12.998	244	1453136	66.366	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	12.627	202	78532	2.459	ng	99
78) Pyrene	12.857	202	79609	2.305	ng	98
83) Chrysene	14.080	228	55100	2.072	ng	99

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

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