

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101725\
 Data File : BF143992.D
 Acq On : 17 Oct 2025 18:43
 Operator : RC/JU
 Sample : Q3349-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PURS Station

Quant Time: Oct 20 11:12:18 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100625.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 06 15:29:47 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	87692	20.000	ng	0.00
21) Naphthalene-d8	8.163	136	299344	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	127355	20.000	ng	0.00
64) Phenanthrene-d10	11.422	188	159085	20.000	ng	0.00
76) Chrysene-d12	14.092	240	119082	20.000	ng	# 0.02
86) Perylene-d12	15.598	264	176056	20.000	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	97093	17.582	ng	0.00
7) Phenol-d6	6.498	99	112124	17.034	ng	-0.02
23) Nitrobenzene-d5	7.440	82	58666	11.907	ng	-0.02
42) 2,4,6-Tribromophenol	10.722	330	16986	13.391	ng	0.00
45) 2-Fluorobiphenyl	9.239	172	106480	13.267	ng	-0.01
79) Terphenyl-d14	13.022	244	91163	13.084	ng	0.01
Target Compounds						
22) Acetophenone	7.287	105	35141	5.552	ng	# 90

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

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