

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102225\
 Data File : BF144037.D
 Acq On : 22 Oct 2025 12:55
 Operator : RC/JU
 Sample : Q3393-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PR132-S06-000021-20251017

Quant Time: Oct 22 13:25:33 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	86146	20.000	ng	0.00
21) Naphthalene-d8	8.163	136	309033	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	160156	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	262433	20.000	ng	0.00
76) Chrysene-d12	14.057	240	218012	20.000	ng	-0.01
86) Perylene-d12	15.545	264	261186	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	373253	68.804	ng	0.00
7) Phenol-d6	6.504	99	416805	64.459	ng	-0.01
23) Nitrobenzene-d5	7.440	82	229915	45.200	ng	-0.02
42) 2,4,6-Tribromophenol	10.710	330	81305	50.969	ng	-0.01
45) 2-Fluorobiphenyl	9.239	172	387848	38.426	ng	-0.01
79) Terphenyl-d14	12.998	244	388948	30.491	ng	-0.01

Target Compounds	Qvalue

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

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