

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060225\
 Data File : BF142600.D
 Acq On : 02 Jun 2025 17:11
 Operator : RC/JU
 Sample : Q2149-01DL 5X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FILTER-CAKEDL

Quant Time: Jun 02 19:03:21 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 20 16:26:47 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.898	152	108020	20.000	ng	0.00
21) Naphthalene-d8	8.181	136	403340	20.000	ng	0.00
39) Acenaphthene-d10	9.939	164	202770	20.000	ng	0.00
64) Phenanthrene-d10	11.427	188	287091	20.000	ng	0.00
76) Chrysene-d12	14.068	240	219560	20.000	ng	0.00
86) Perylene-d12	15.568	264	244032	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	166002	25.891	ng	0.00
7) Phenol-d6	6.522	99	181610	23.531	ng	-0.01
23) Nitrobenzene-d5	7.457	82	138869	18.774	ng	-0.02
42) 2,4,6-Tribromophenol	10.727	330	54718	24.314	ng	-0.01
45) 2-Fluorobiphenyl	9.251	172	287136	19.002	ng	-0.02
79) Terphenyl-d14	13.010	244	188080	11.706	ng	0.00
Target Compounds						
15) Benzyl Alcohol	7.034	79	19655	3.446	ng	Qvalue 97
20) 3+4-Methylphenols	7.298	107	236135	33.819	ng	# 54
32) Benzoic acid	7.934	122	10971	2.867	ng	98

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

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