

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102125\
 Data File : BF144014.D
 Acq On : 21 Oct 2025 16:03
 Operator : RC/JU
 Sample : Q3386-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 3444

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 10/22/2025
 Supervised By :Jagrut Upadhyay 10/22/2025

Quant Time: Oct 21 16:30:07 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	109277	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	380104	20.000	ng	0.00
39) Acenaphthene-d10	9.928	164	175218	20.000	ng	0.00
64) Phenanthrene-d10	11.422	188	254880	20.000	ng	0.00
76) Chrysene-d12	14.086	240	175156	20.000	ng	# 0.02
86) Perylene-d12	15.586	264	269690	20.000	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	587367	85.355	ng	0.00
7) Phenol-d6	6.516	99	578488m	70.526	ng	0.00
23) Nitrobenzene-d5	7.440	82	441754	70.609	ng	-0.02
42) 2,4,6-Tribromophenol	10.722	330	185649	106.377	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	819469	74.210	ng	0.00
79) Terphenyl-d14	13.016	244	714511	69.718	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.534	94	32481m	3.312	ng	
15) Benzyl Alcohol	7.034	79	41472	6.748	ng	94
32) Benzoic acid	8.157	122	187803m	50.643	ng	
70) Pentachlorophenol	11.228	266	6743	3.600	ng	94
84) Bis(2-ethylhexyl)phtha...	14.057	149	162000	28.338	ng	# 89

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

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