

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102125\
 Data File : BF144013.D
 Acq On : 21 Oct 2025 15:33
 Operator : RC/JU
 Sample : Q3386-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 3443

Manual Integrations
 APPROVED

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

Quant Time: Oct 21 16:22:07 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	84784	20.000	ng	0.00
21) Naphthalene-d8	8.163	136	313703	20.000	ng	0.00
39) Acenaphthene-d10	9.928	164	132876	20.000	ng	0.00
64) Phenanthrene-d10	11.422	188	197689	20.000	ng	0.00
76) Chrysene-d12	14.092	240	140758	20.000	ng	# 0.02
86) Perylene-d12	15.598	264	234614	20.000	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	461517	86.441	ng	0.00
7) Phenol-d6	6.510	99	459042	72.132	ng	0.00
23) Nitrobenzene-d5	7.439	82	340852	66.013	ng	-0.02
42) 2,4,6-Tribromophenol	10.722	330	161919	122.344	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	664467	79.347	ng	0.00
79) Terphenyl-d14	13.021	244	620483	75.339	ng	0.01
Target Compounds						
10) Phenol	6.528	94	33467	4.399	ng	# 40
15) Benzyl Alcohol	7.028	79	33546m	7.035	ng	
32) Benzoic acid	8.110	122	204964m	66.970	ng	
70) Pentachlorophenol	11.227	266	4759	3.276	ng	96
84) Bis(2-ethylhexyl)phtha...	14.063	149	111450	24.260	ng	# 80

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

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