

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091825\
 Data File : BP025781.D
 Acq On : 18 Sep 2025 13:03
 Operator : CG/JU
 Sample : Q3015-06
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 WP0925-PT-ACIDS-WP

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP025781.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.990	4	9	14	rBV	1929125	2325801	23.33%	2.415%
2	3.031	14	16	32	rVB3	99966	149927	1.50%	0.156%
3	4.872	322	329	343	rBV	424298	655087	6.57%	0.680%
4	5.325	399	406	420	rBV	3014443	4337824	43.51%	4.505%
5	6.890	662	672	674	rBV	2554856	4331552	43.45%	4.498%
6	7.278	731	738	755	rBV	1014804	1740128	17.46%	1.807%
7	7.702	802	810	818	rBV	503026	812683	8.15%	0.844%
8	8.143	878	885	905	rBV	1124795	1826950	18.33%	1.897%
9	8.460	932	939	965	rBV	1639815	2892129	29.01%	3.003%
10	8.849	997	1005	1013	rBV	1644677	3055083	30.65%	3.173%
11	9.596	1125	1132	1138	rBV	2013362	3336249	33.47%	3.465%
12	9.666	1138	1144	1167	rVB	2650050	4915764	49.31%	5.105%
13	10.143	1217	1225	1244	rBV	696375	1422765	14.27%	1.477%
14	10.484	1274	1283	1291	rBV	662656	1100355	11.04%	1.143%
15	10.666	1303	1314	1331	rBV	2880339	6108022	61.27%	6.343%
16	12.778	1665	1673	1681	rBV	4182183	7058162	70.80%	7.330%
17	12.860	1681	1687	1698	rVV	3038448	5567149	55.84%	5.781%
18	12.960	1698	1704	1714	rVB	4493428	6466936	64.87%	6.716%
19	13.678	1820	1826	1835	rBV	66531	116893	1.17%	0.121%
20	14.360	1935	1942	1949	rBV	1039822	1426289	14.31%	1.481%
21	14.478	1956	1962	1973	rBV	1803102	3178600	31.88%	3.301%
22	14.613	1979	1985	1997	rBV	1149759	2661497	26.70%	2.764%
23	15.007	2045	2052	2071	rBV	3206006	5643057	56.61%	5.860%
24	15.519	2132	2139	2156	rBV	1139701	1831159	18.37%	1.902%
25	15.878	2192	2200	2212	rBV	3101489	5024772	50.40%	5.218%
26	16.825	2354	2361	2380	rBV	1577591	2542425	25.50%	2.640%
27	17.178	2414	2421	2434	rBV2	1070184	1733029	17.38%	1.800%
28	19.913	2879	2886	2895	rBV	6925438	9968995	100.00%	10.352%
29	21.642	3174	3180	3196	rVB	1107624	2139236	21.46%	2.222%
30	25.012	3744	3753	3768	rVB2	652754	1927314	19.33%	2.001%

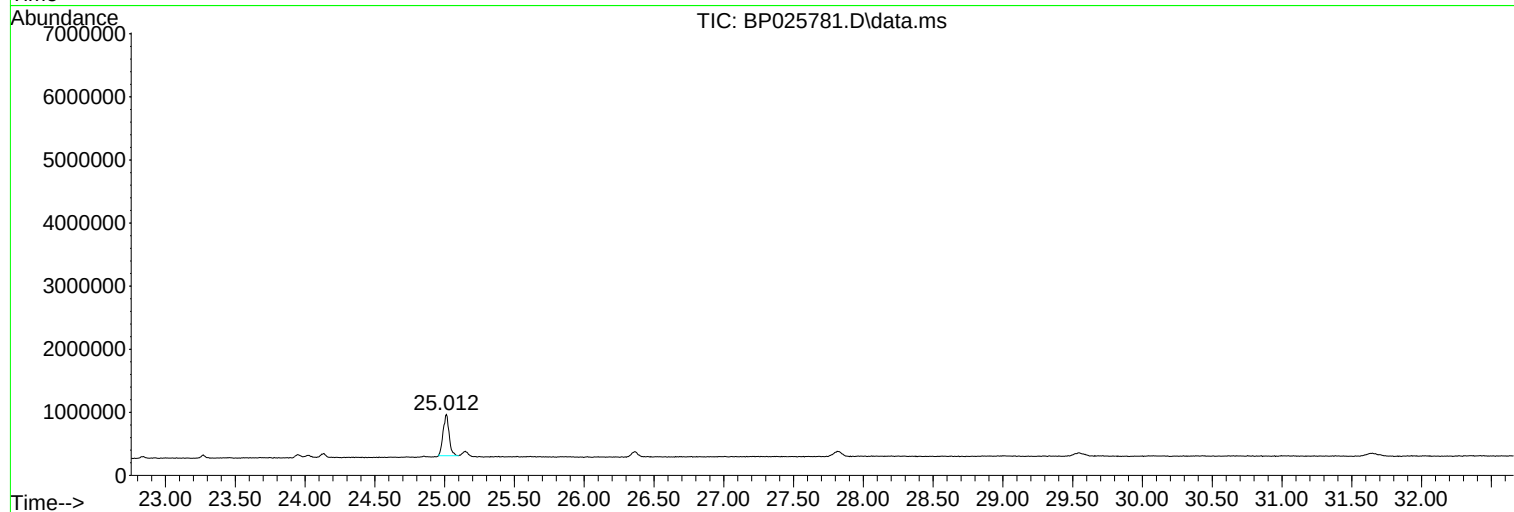
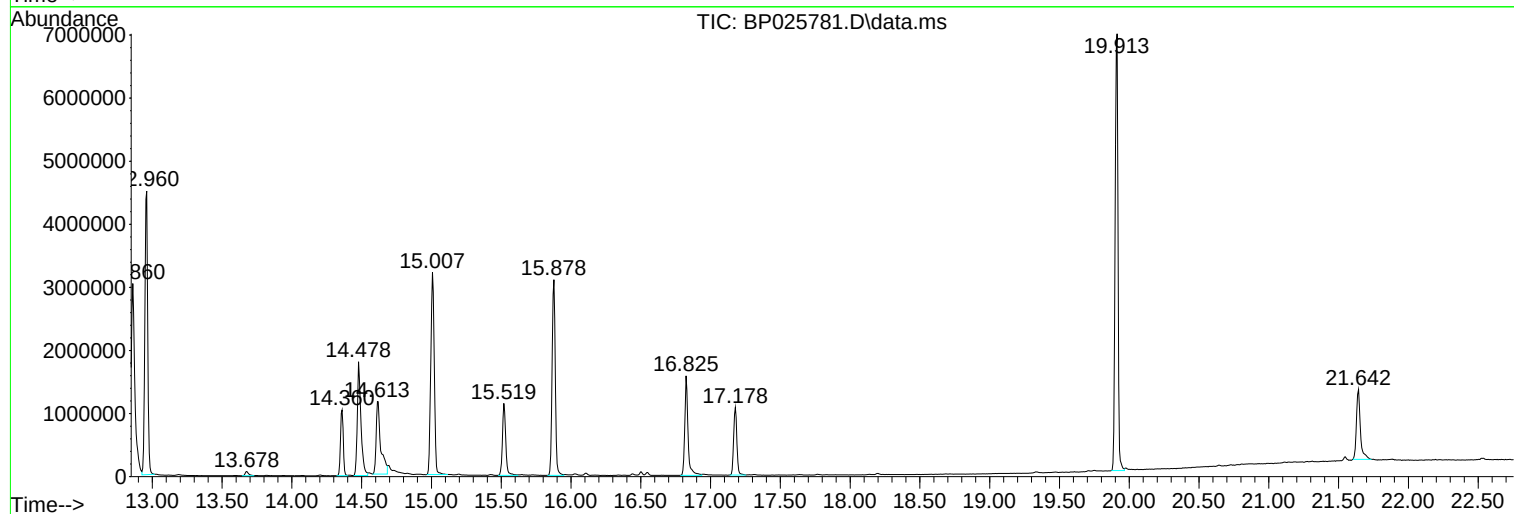
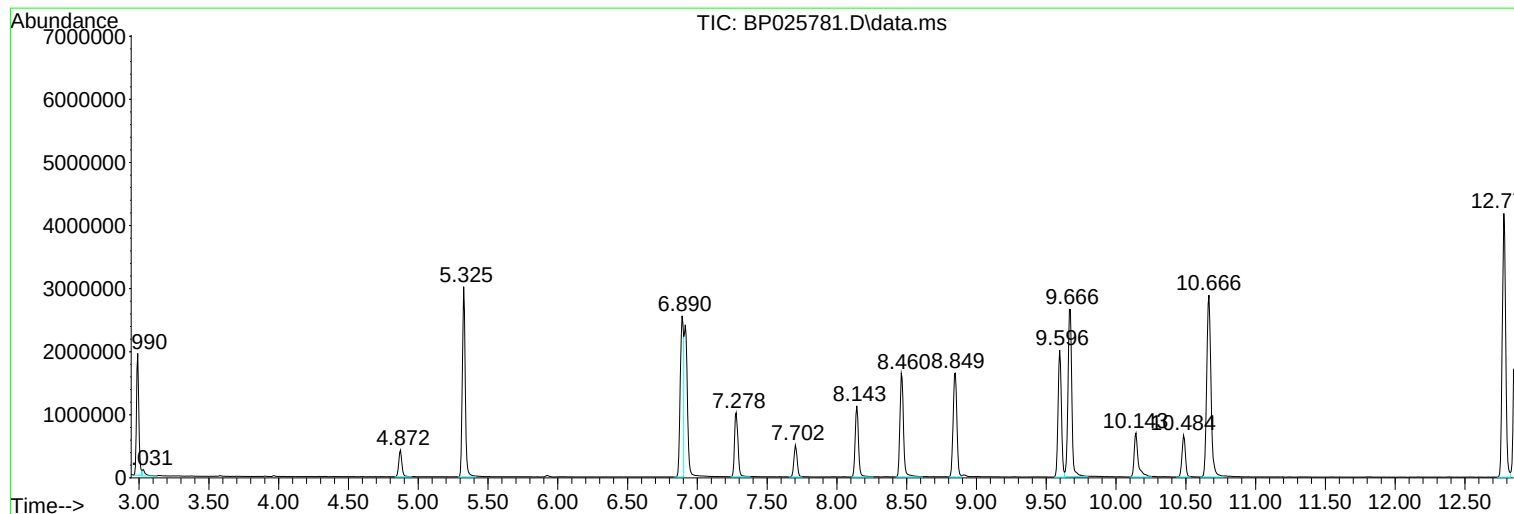
Sum of corrected areas: 96295832

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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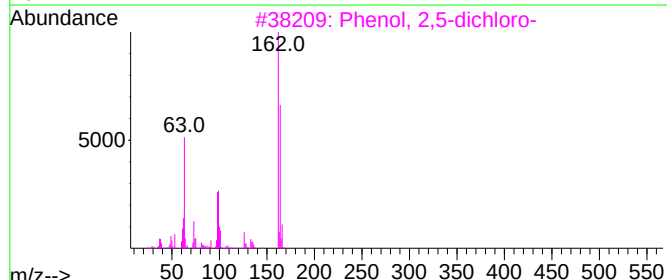
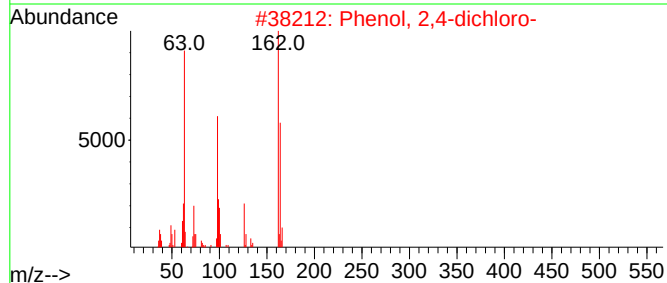
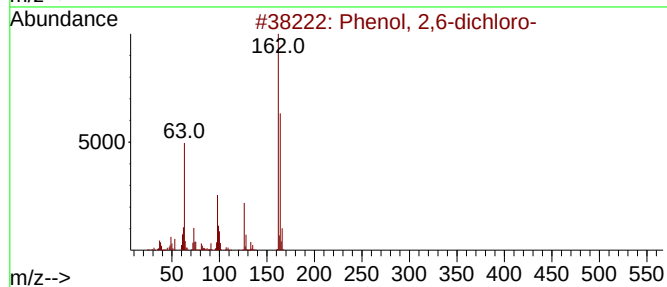
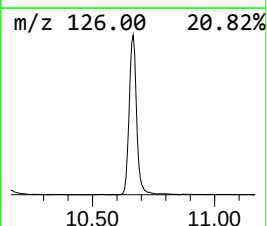
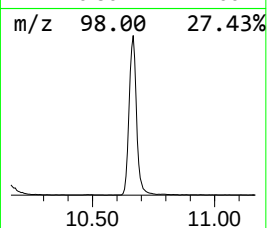
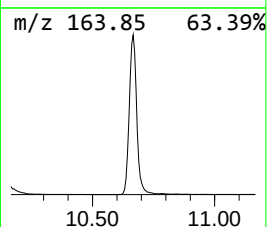
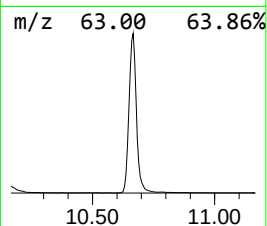
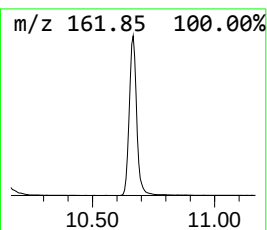
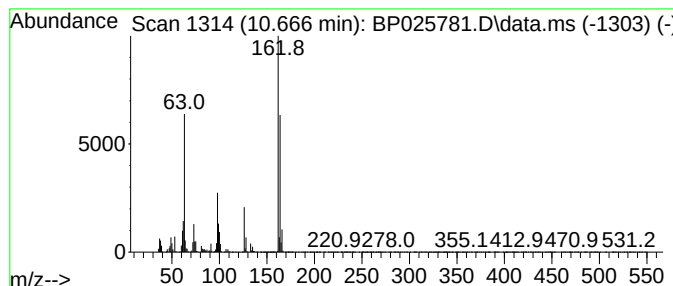
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Phenol, 2,6-dichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.666	111.02 ng	6108020	Naphthalene-d8	10.484

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,6-dichloro-	162	C6H4Cl2O	000087-65-0	98
2			Phenol, 2,4-dichloro-	162	C6H4Cl2O	000120-83-2	97
3			Phenol, 2,5-dichloro-	162	C6H4Cl2O	000583-78-8	94
4			Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	94
5			Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	50



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenol, 2,6-dic...	10.666	111.0	ng	6108020	2	10.484	1100360	20.0