

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080624\
 Data File : BP021398.D
 Acq On : 06 Aug 2024 20:07
 Operator : CG/JU
 Sample : P3359-17
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E1GF6

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 08/07/2024
 Supervised By :mohammad ahmed 08/09/2024

Quant Time: Aug 06 23:23:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 29 12:46:18 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	197852	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.381	136	805637	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.263	164	543754	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.086	188	1220340	20.000	ng/u1	0.00
79) Chrysene-d12	21.521	240	1325041	20.000	ng/u1	-0.01
88) Perylene-d12	24.815	264	1565177	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.152	96	19832	4.561	ng/uL	-0.02
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	6.852	99	185578	11.505	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	270479	27.669	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.169	132	370525	27.880	ng/u1	-0.01
15) 4-Methylphenol-d8	8.363	113	297699	22.221	ng/u1	0.00
21) Nitrobenzene-d5	8.781	128	175211	28.000	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.499	143	208905	29.168	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.069	165	353100m	27.265	ng/u1	0.02
31) 4-Chloroaniline-d4	10.599	131	47737m	2.782	ng/u1	0.04
46) Dimethylphthalate-d6	13.681	166	1135696	28.731	ng/u1	0.00
49) Acenaphthylene-d8	13.957	160	1200593	27.050	ng/u1	0.00
54) 4-Nitrophenol-d4	14.675	143	67981m	12.088	ng/u1	0.09
60) Fluorene-d10	15.281	176	983542	30.330	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.481	200	169132m	22.906	ng/u1	0.02
73) Anthracene-d10	17.186	188	1580289	29.156	ng/u1	0.00
81) Pyrene-d10	19.569	212	2104802	25.694	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.568	264	2304563	29.854	ng/u1	-0.03
Target Compounds						
6) Benzaldehyde	6.816	77	12868	1.582	ng/u1	91
16) Acetophenone	8.463	105	21017	1.016	ng/u1	95

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

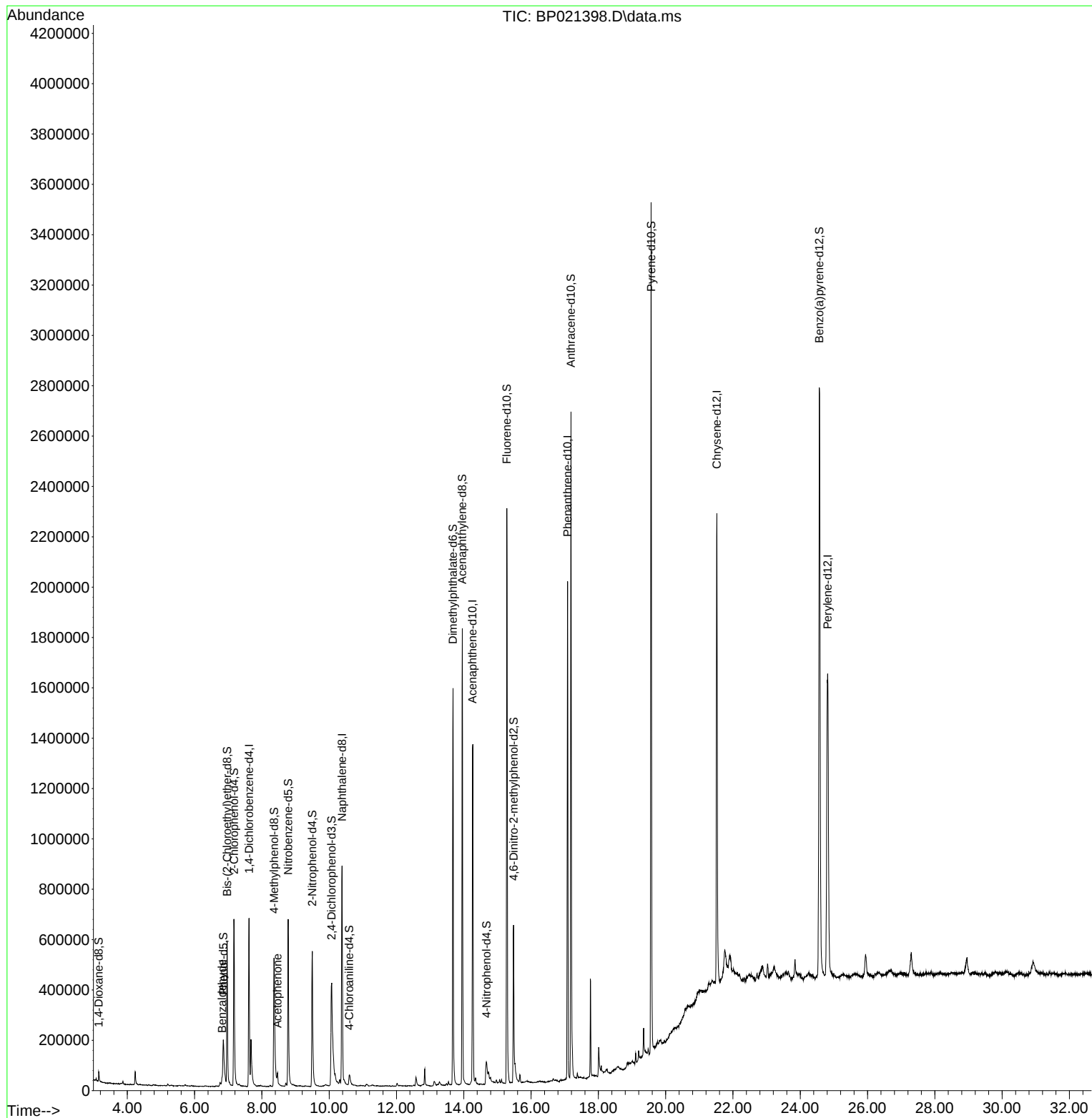
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080624\
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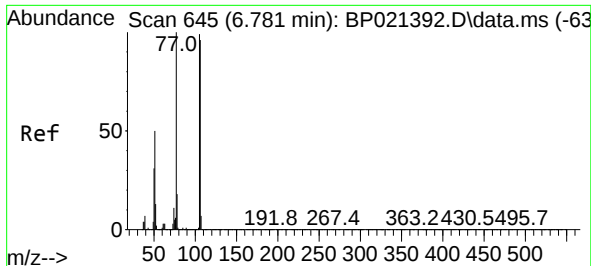
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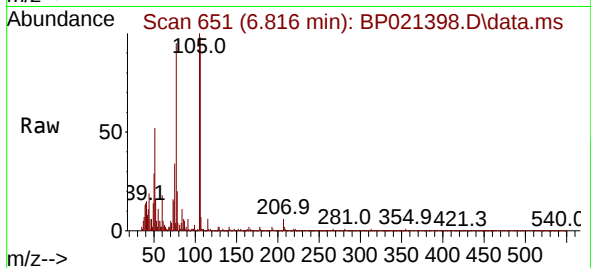
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Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 29 12:46:18 2024
Response via : Initial Calibration





#6
Benzaldehyde
Concen: 1.582 ng/ul
RT: 6.816 min Scan# 61
Delta R.T. 0.012 min
Lab File: BP021398.D
Acq: 06 Aug 2024 20:07

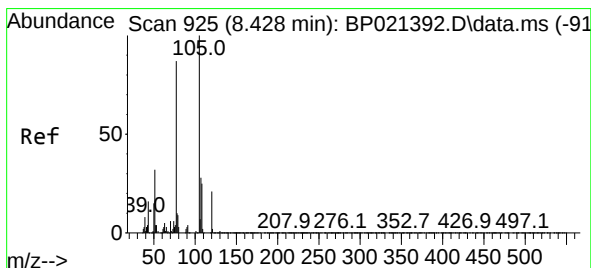
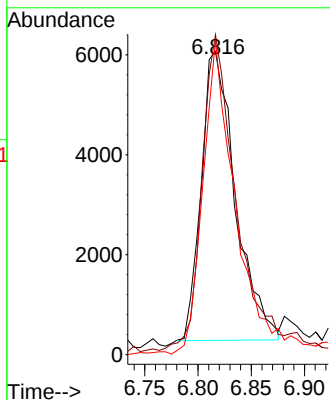
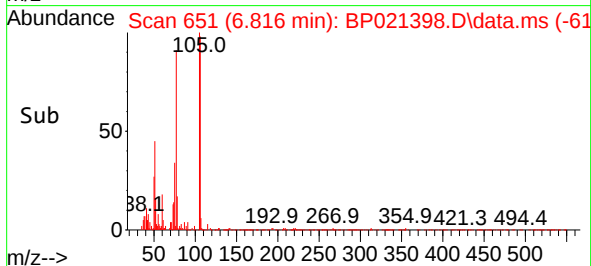
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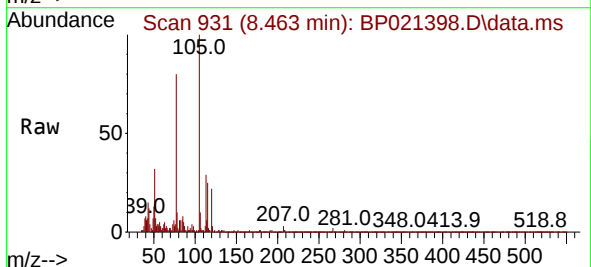
Tgt Ion: 77 Resp: 1286
Ion Ratio Lower Upper
77 100
105 105.6 79.3 118.9
106 103.7 74.6 112.0

Manual Integrations
APPROVED

Reviewed By :Yogesh Patel 08/07/2024
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#16
Acetophenone
Concen: 1.016 ng/ul
RT: 8.463 min Scan# 931
Delta R.T. 0.006 min
Lab File: BP021398.D
Acq: 06 Aug 2024 20:07



Tgt Ion:105 Resp: 21017
Ion Ratio Lower Upper
105 100
77 79.7 61.8 92.8
51 32.0 21.5 32.3

