

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
 Data File : BP019048.D
 Acq On : 22 Dec 2023 20:01
 Operator : MA/JU
 Sample : 05835-13
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AR0

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 12/25/2023
 Supervised By :mohammad ahmed 12/25/2023

Quant Time: Dec 22 22:03:59 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 21 04:46:12 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	149786	20.000	ng/u1	0.00
20) Naphthalene-d8	10.610	136	573017	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.451	164	377008	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.204	188	890727m	20.000	ng/u1	0.00
79) Chrysene-d12	21.292	240	921447	20.000	ng/u1	0.00
88) Perylene-d12	23.651	264	1060106	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	12370	2.814	ng/uL	0.00
4) Pyridine-d5	3.675	84	56162	4.770	ng/u1	0.00
7) Phenol-d5	6.993	99	221351	14.726	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.157	67	134179	15.061	ng/u1	0.00
11) 2-Chlorophenol-d4	7.352	132	185576	15.822	ng/u1	0.00
15) 4-Methylphenol-d8	8.534	113	146301	12.027	ng/u1	0.00
21) Nitrobenzene-d5	8.981	128	88146	17.283	ng/u1	0.00
24) 2-Nitrophenol-d4	9.699	143	98567	18.641	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.240	165	191190	17.683	ng/u1	0.00
31) 4-Chloroaniline-d4	10.763	131	68429	4.843	ng/u1	0.00
46) Dimethylphthalate-d6	13.869	166	599561	17.818	ng/u1	0.00
49) Acenaphthylene-d8	14.145	160	666835	17.998	ng/u1	0.00
54) 4-Nitrophenol-d4	14.675	143	69536	12.590	ng/u1	0.00
60) Fluorene-d10	15.445	176	543324	19.237	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	75984m	14.227	ng/u1	0.00
73) Anthracene-d10	17.298	188	818273	17.474	ng/u1	0.00
81) Pyrene-d10	19.539	212	1078618	15.038	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.498	264	984941	15.935	ng/u1	0.00
Target Compounds					Qvalue	
6) Benzaldehyde	6.963	77	11370	1.465	ng/u1	93
8) Phenol	7.022	94	47369	3.213	ng/u1	97
16) Acetophenone	8.646	105	245966	13.567	ng/u1	99

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

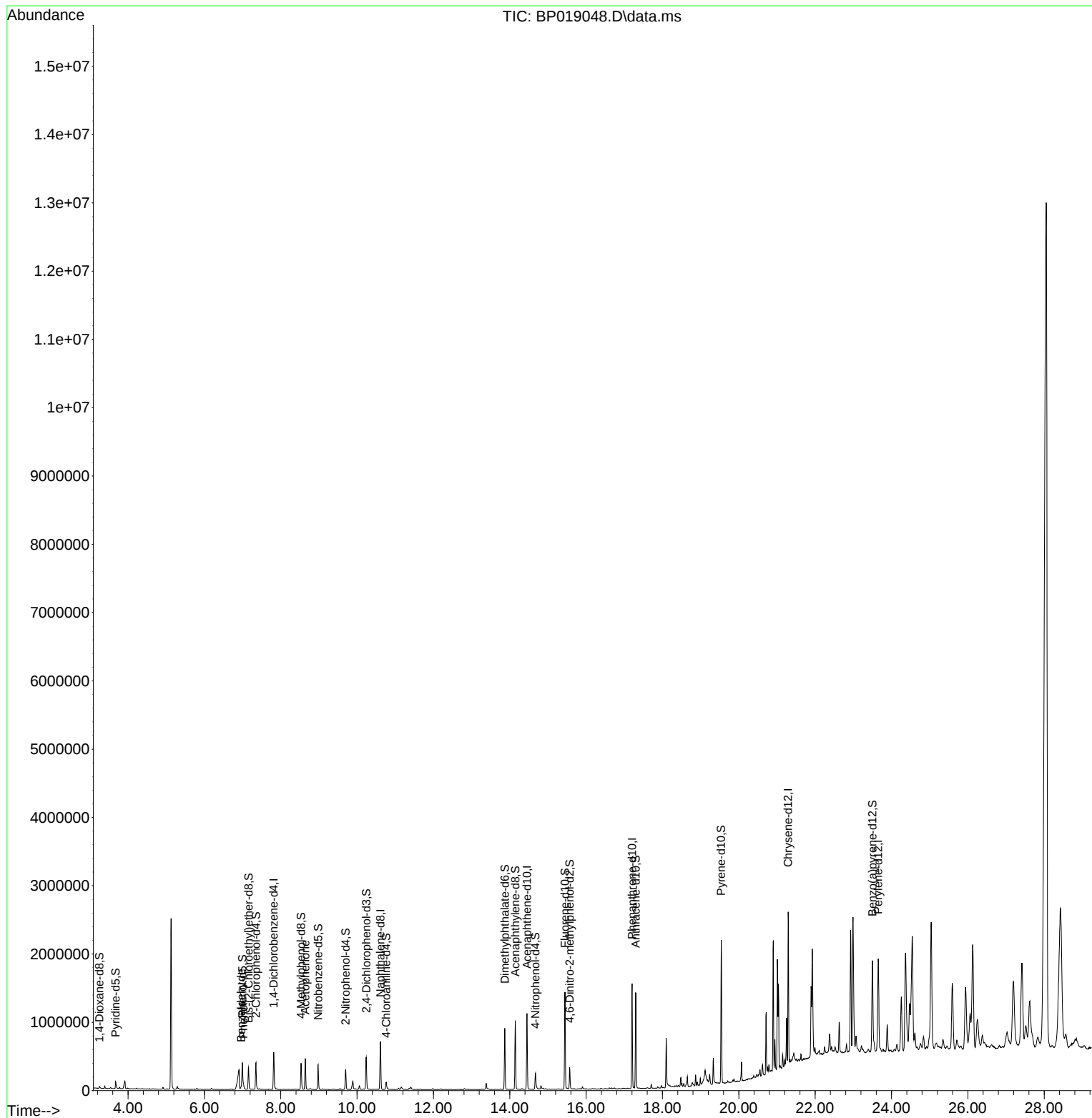
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019048.D
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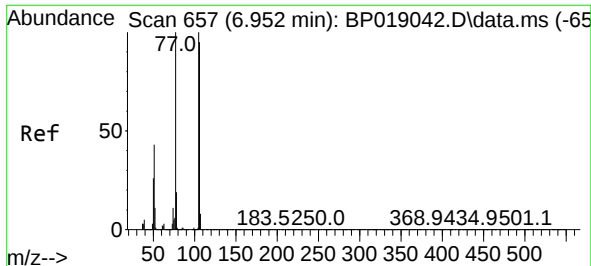
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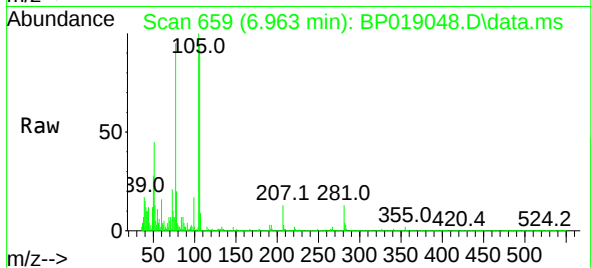
Quant Time: Dec 22 22:03:59 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 21 04:46:12 2023
Response via : Initial Calibration





#6
Benzaldehyde
Concen: 1.465 ng/ul
RT: 6.963 min Scan# 613
Delta R.T. -0.000 min
Lab File: BP019048.D
Acq: 22 Dec 2023 20:01

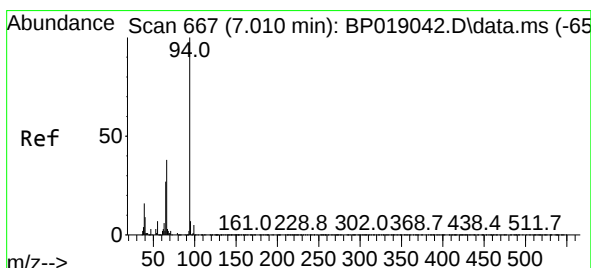
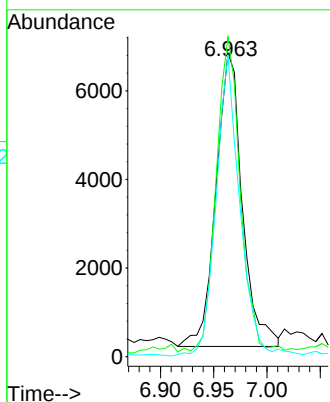
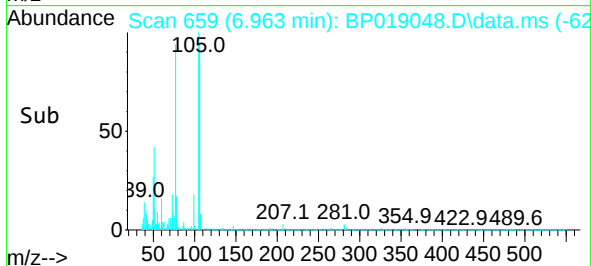
Instrument :
BNA_P
ClientSampleId :
C0AR0



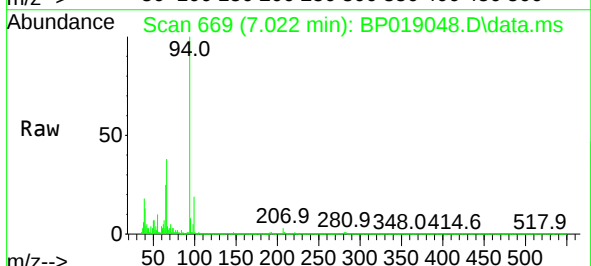
Tgt Ion: 77 Resp: 11370
Ion Ratio Lower Upper
77 100
105 105.3 77.9 116.9
106 99.4 74.6 111.8

Manual Integrations
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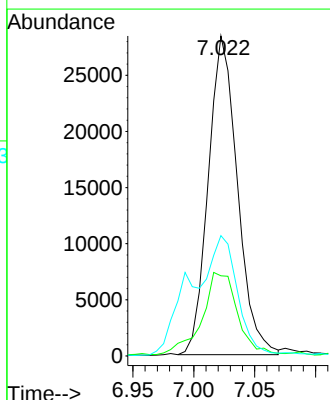
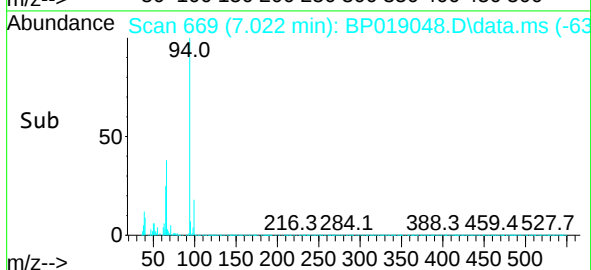
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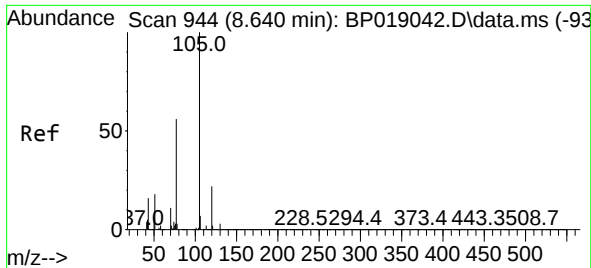


#8
Phenol
Concen: 3.213 ng/ul
RT: 7.022 min Scan# 669
Delta R.T. -0.006 min
Lab File: BP019048.D
Acq: 22 Dec 2023 20:01



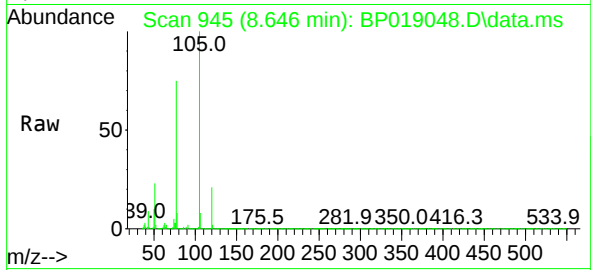
Tgt Ion: 94 Resp: 47369
Ion Ratio Lower Upper
94 100
65 25.0 21.0 31.6
66 37.6 28.6 43.0





#16
Acetophenone
Concen: 13.567 ng/ul
RT: 8.646 min Scan# 944
Delta R.T. -0.006 min
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Tgt Ion:105 Resp: 245960
Ion Ratio Lower Upper
105 100
77 74.9 59.4 89.0
51 23.4 19.4 29.0

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