

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091823\
 Data File : BP017228.D
 Acq On : 19 Sep 2023 04:14
 Operator : MA/JU
 Sample : 04332-10
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AH2

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 09/19/2023
 Supervised By :Sohil Jodhani 09/19/2023

Quant Time: Sep 19 05:08:44 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 16 03:18:48 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	237012	20.000	ng/u1	0.00
20) Naphthalene-d8	10.769	136	924731	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.604	164	483845	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.375	188	913678	20.000	ng/u1	0.00
79) Chrysene-d12	21.474	240	858234	20.000	ng/u1	0.00
88) Perylene-d12	24.004	264	1052251	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	9042	1.603	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	7.099	99	131666	6.832	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	106148	8.812	ng/u1	0.00
11) 2-Chlorophenol-d4	7.463	132	125635	8.172	ng/u1	0.00
15) 4-Methylphenol-d8	8.657	113	66131	4.140	ng/u1	0.00
21) Nitrobenzene-d5	9.140	128	62821	9.702	ng/u1	0.00
24) 2-Nitrophenol-d4	9.857	143	65097	9.417	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	101910	7.623	ng/u1	0.00
31) 4-Chloroaniline-d4	10.940	131	32084	1.592	ng/u1	0.00
46) Dimethylphthalate-d6	14.016	166	321851	9.348	ng/u1	0.00
49) Acenaphthylene-d8	14.304	160	351458	8.996	ng/u1	0.00
54) 4-Nitrophenol-d4	14.839	143	21186m	3.184	ng/u1	0.01
60) Fluorene-d10	15.598	176	259686	8.723	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.733	200	15266	2.991	ng/u1	0.00
73) Anthracene-d10	17.475	188	334422	8.615	ng/u1	0.00
81) Pyrene-d10	19.716	212	428509	9.520	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.833	264	444207	8.944	ng/u1	0.00
Target Compounds						
					Qvalue	
6) Benzaldehyde	7.105	77	57485	5.358	ng/u1	99
16) Acetophenone	8.793	105	59986	2.399	ng/u1	96

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

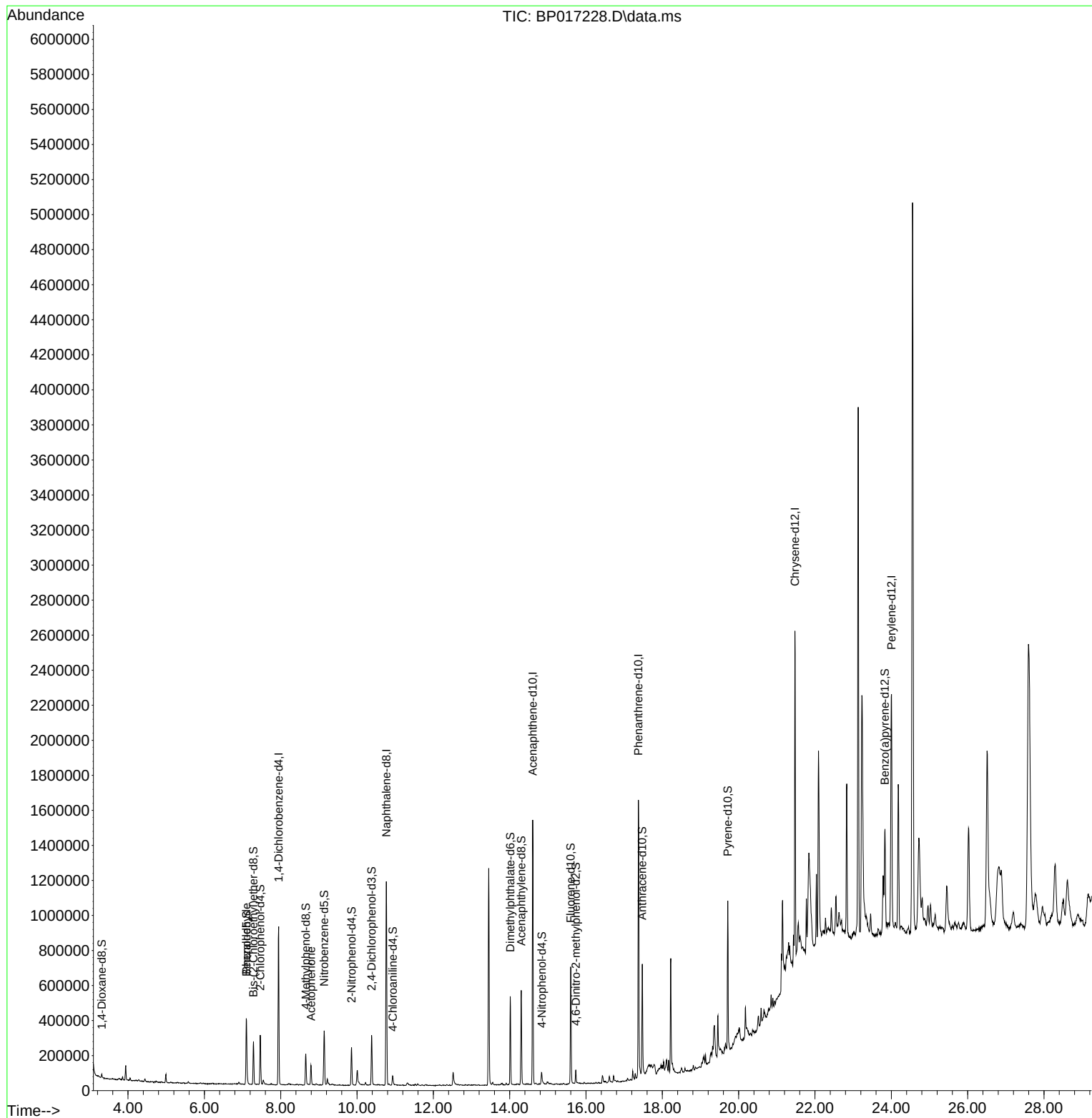
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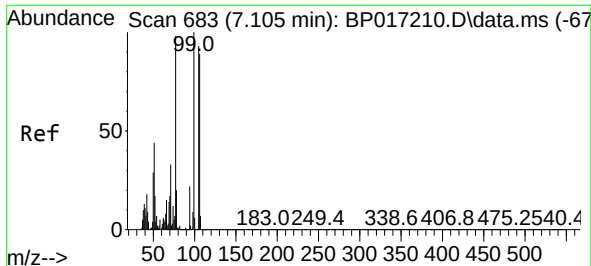
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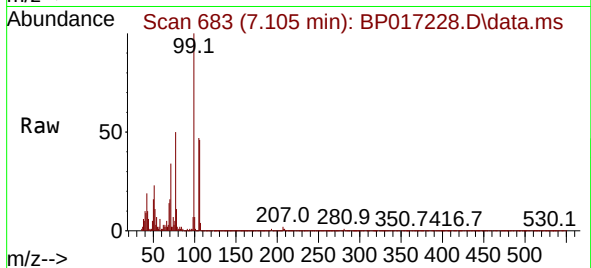
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Quant Title : SVOA CALIBRATION
QLast Update : Sat Sep 16 03:18:48 2023
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#6
Benzaldehyde
Concen: 5.358 ng/ul
RT: 7.105 min Scan# 683
Delta R.T. -0.000 min
Lab File: BP017228.D
Acq: 19 Sep 2023 04:14

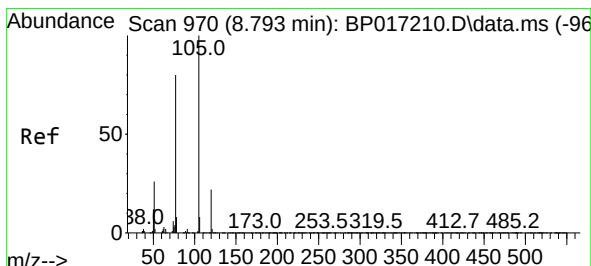
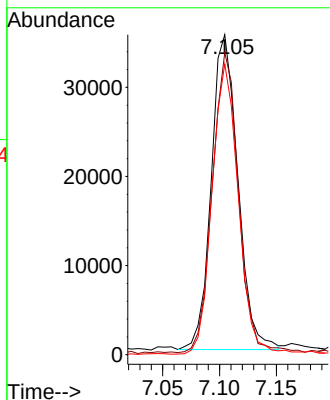
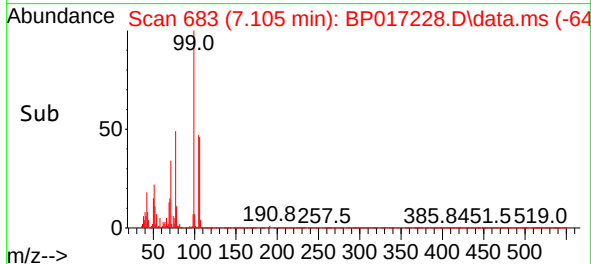
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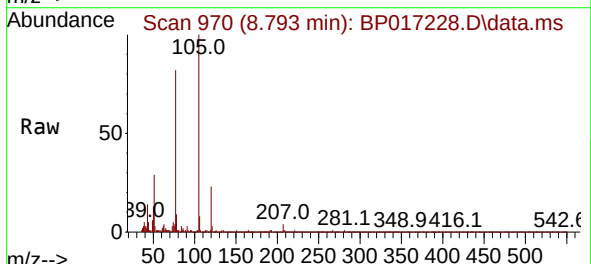
Tgt Ion: 77 Resp: 57485
Ion Ratio Lower Upper
77 100
105 94.4 75.3 112.9
106 90.9 74.6 111.8

Manual Integrations
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#16
Acetophenone
Concen: 2.399 ng/ul
RT: 8.793 min Scan# 970
Delta R.T. -0.006 min
Lab File: BP017228.D
Acq: 19 Sep 2023 04:14



Tgt Ion:105 Resp: 59986
Ion Ratio Lower Upper
105 100
77 81.7 63.0 94.4
51 29.2 20.5 30.7

