

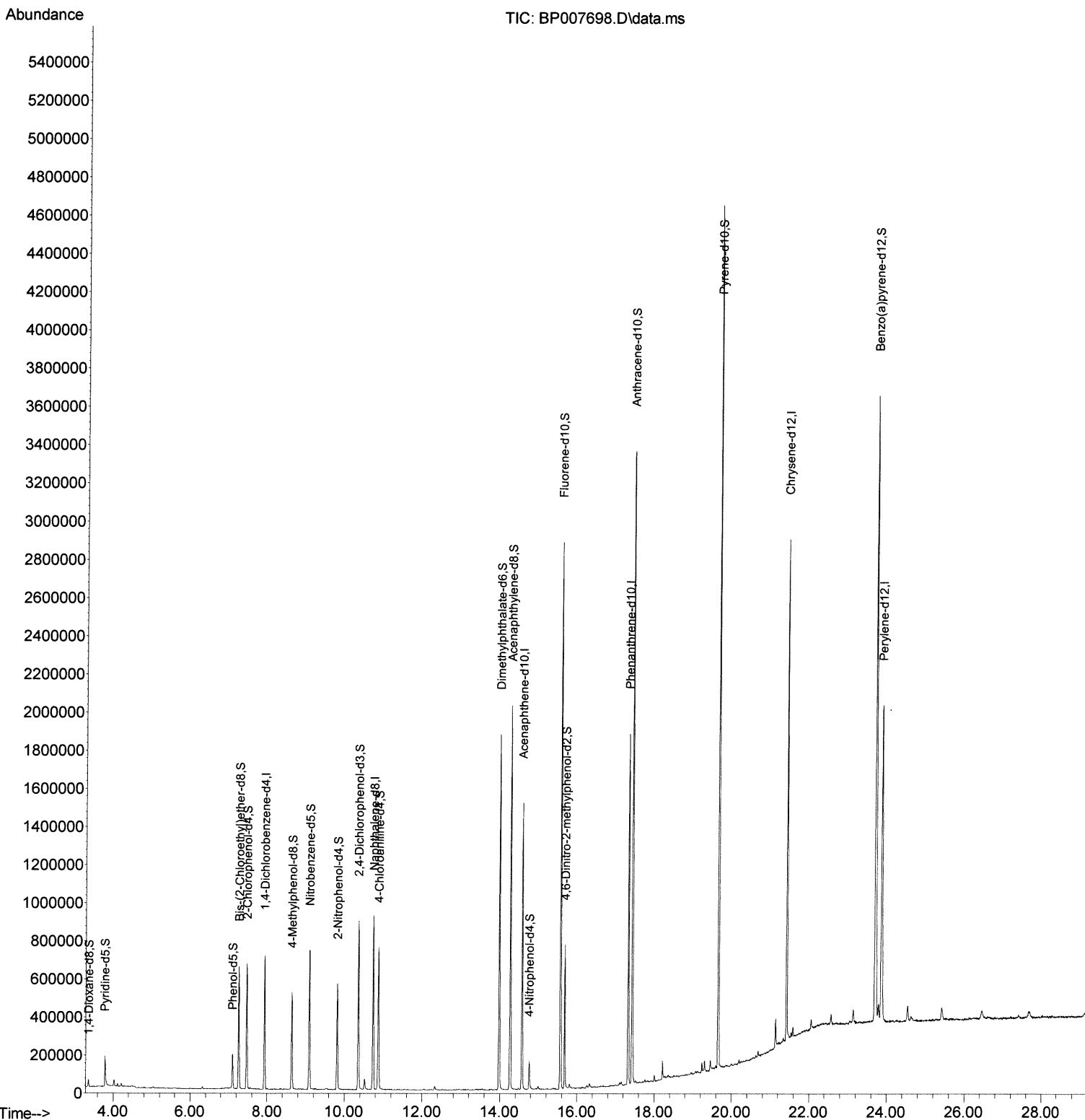
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007698.D
Acq On : 27 Oct 2021 02:24
Operator : CG/JU
Sample : M4303-06
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
P012-TW001-02

Manual Integrations
APPROVED

mohammad
10/27/2021 11:51:03 AM

Quant Time: Oct 27 03:24:22 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Oct 25 16:57:16 2021
Response via : Initial Calibration



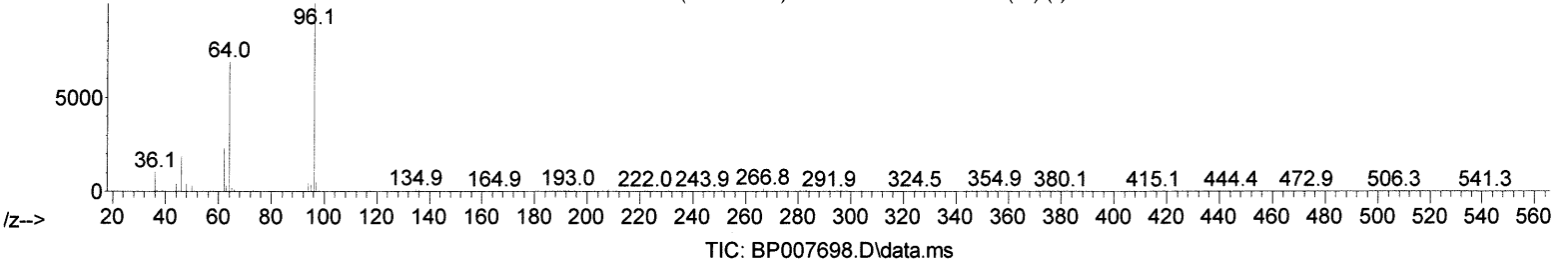
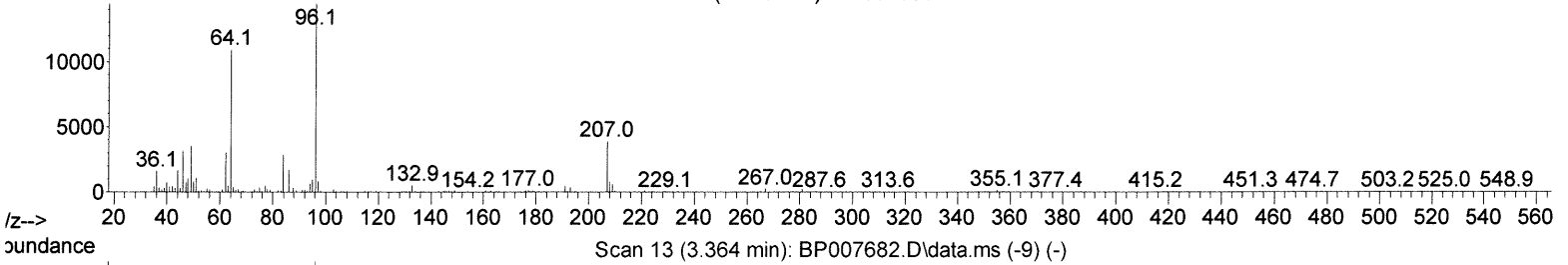
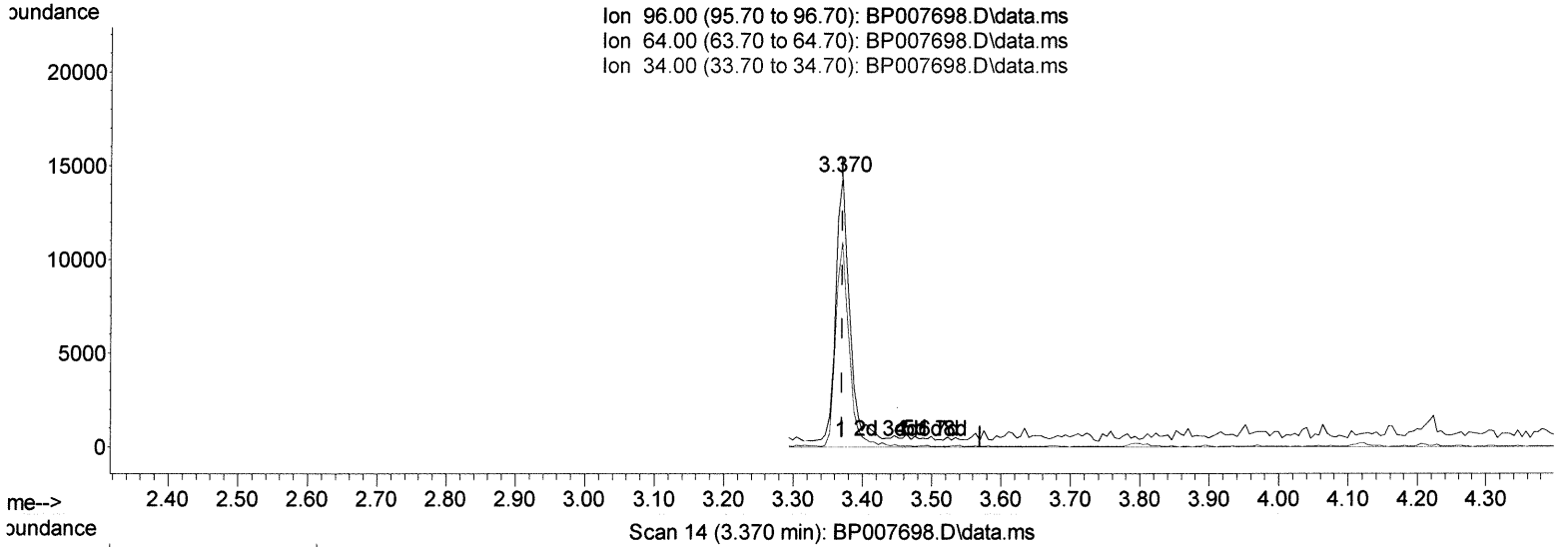
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(3) 1,4-Dioxane-d8 (S)

3.370min (-0.000) 4.41 ng/uL

response 18706

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.20	75.76
34.00	0.00	0.00
0.00	0.00	0.00

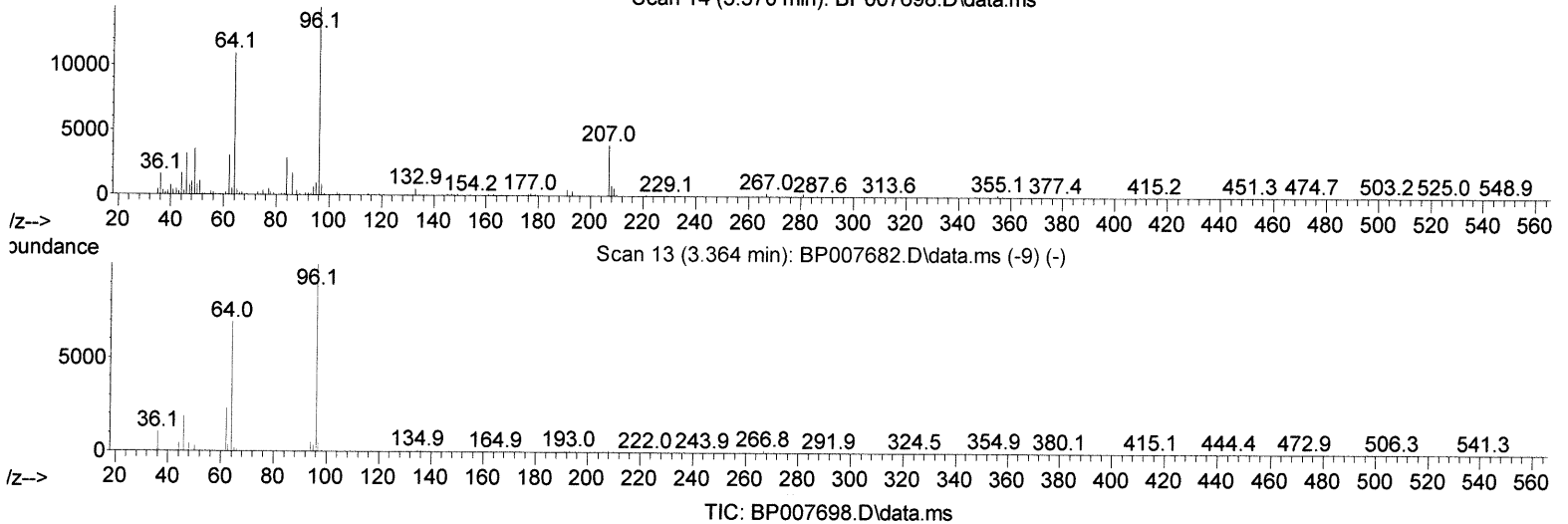
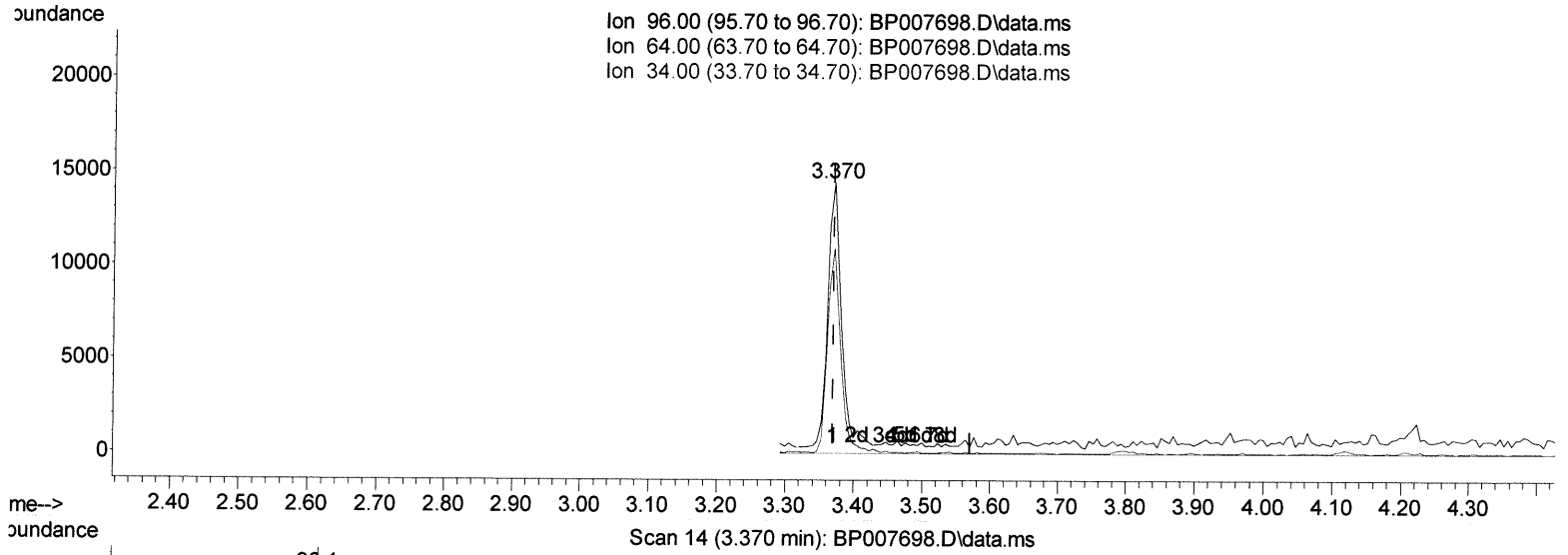
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 Response via : Initial Calibration



(3) 1,4-Dioxane-d8 (S)

3.370min (-0.000) 4.64 ng/uL m 10/29/21 JU

response	19682	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	69.20	75.76
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
 Data File : BP007698.D
 Acq On : 27 Oct 2021 02:24
 Operator : CG/JU
 Sample : M4303-06
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Manual Integrations
 APPROVED

mohammad
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 Quant Title : SVOA CALIBRATION
 Last Update : Mon Oct 25 16:57:16 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	192800	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	774815	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	521247	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.322	188	1144902	20.000	ng/ul	-0.01
79) Chrysene-d12	21.416	240	1251254	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	1296540	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	19682m >	4.637	ng/uL	> 0.00 10/29/21 J4
4) Pyridine-d5	3.787	84	89876	7.002	ng/ul	0.00
7) Phenol-d5	7.099	99	104541	7.139	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	276550	34.014	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	297309	26.103	ng/ul	-0.01
15) 4-Methylphenol-d8	8.640	113	189280	16.507	ng/ul	-0.01
21) Nitrobenzene-d5	9.093	128	180386	34.163	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	179117	31.059	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.357	165	333865	28.442	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.869	131	403946	26.026	ng/ul	-0.01
46) Dimethylphthalate-d6	13.981	166	1233746	34.943	ng/ul	-0.01
49) Acenaphthylene-d8	14.263	160	1433898	33.134	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	39057	6.035	ng/ul	-0.02
60) Fluorene-d10	15.563	176	1115820	37.487	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.681	200	166933	27.258	ng/ul	0.00
73) Anthracene-d10	17.428	188	1973507	41.944	ng/ul	0.00
81) Pyrene-d10	19.657	212	2456833	42.174	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	2517063	39.519	ng/ul	0.00

Target Compounds Qvalue

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