

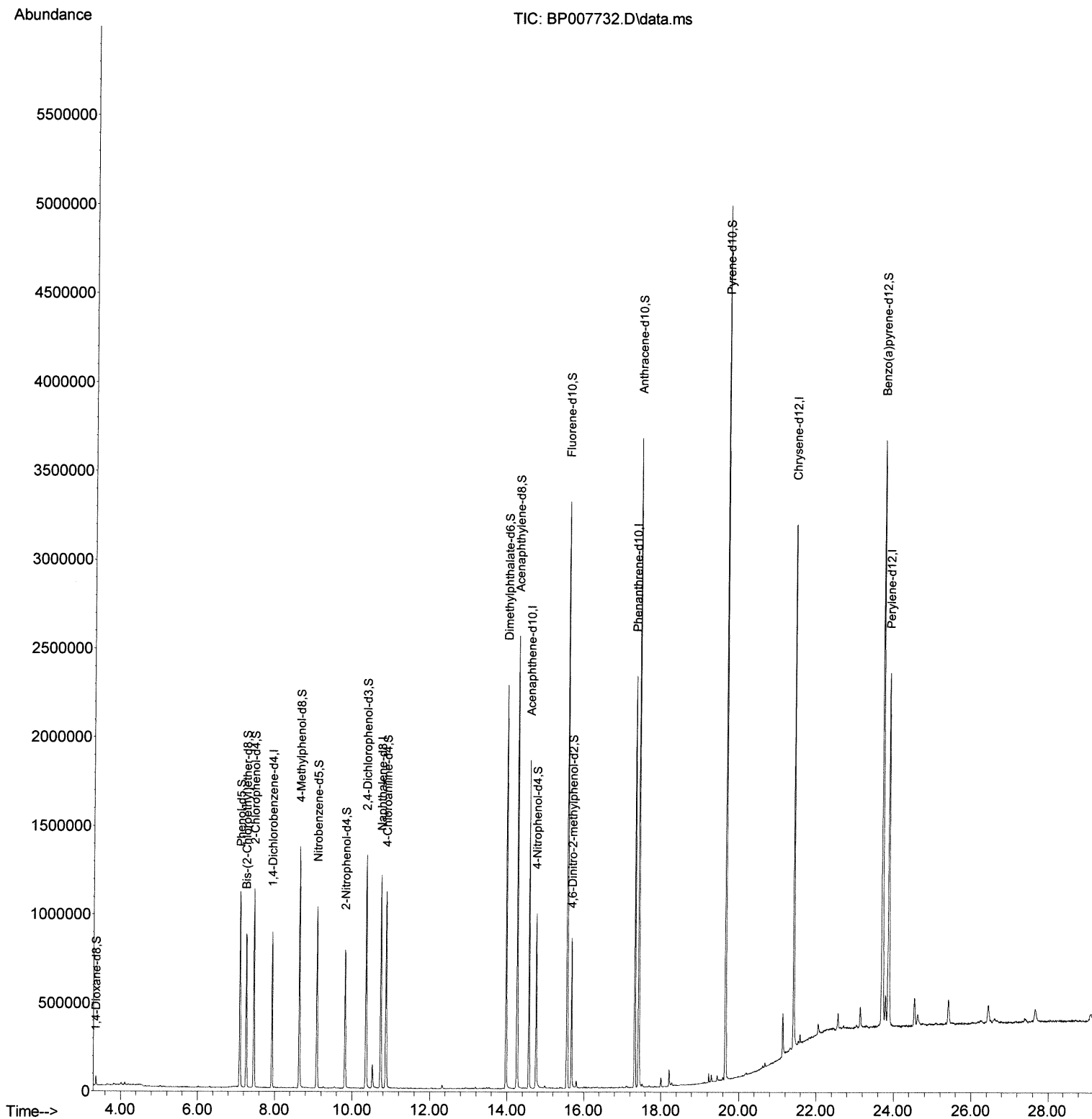
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102521\
Data File : BP007732.D
Acq On : 27 Oct 2021 22:19
Operator : CG/JU
Sample : M4215-18
Misc :
ALS Vial : 89 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JDGF5

Manual Integrations
APPROVED

mohammad
10/28/2021 11:12:18 AM

Quant Time: Oct 28 01:36:50 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 27 15:37:43 2021
Response via : Initial Calibration

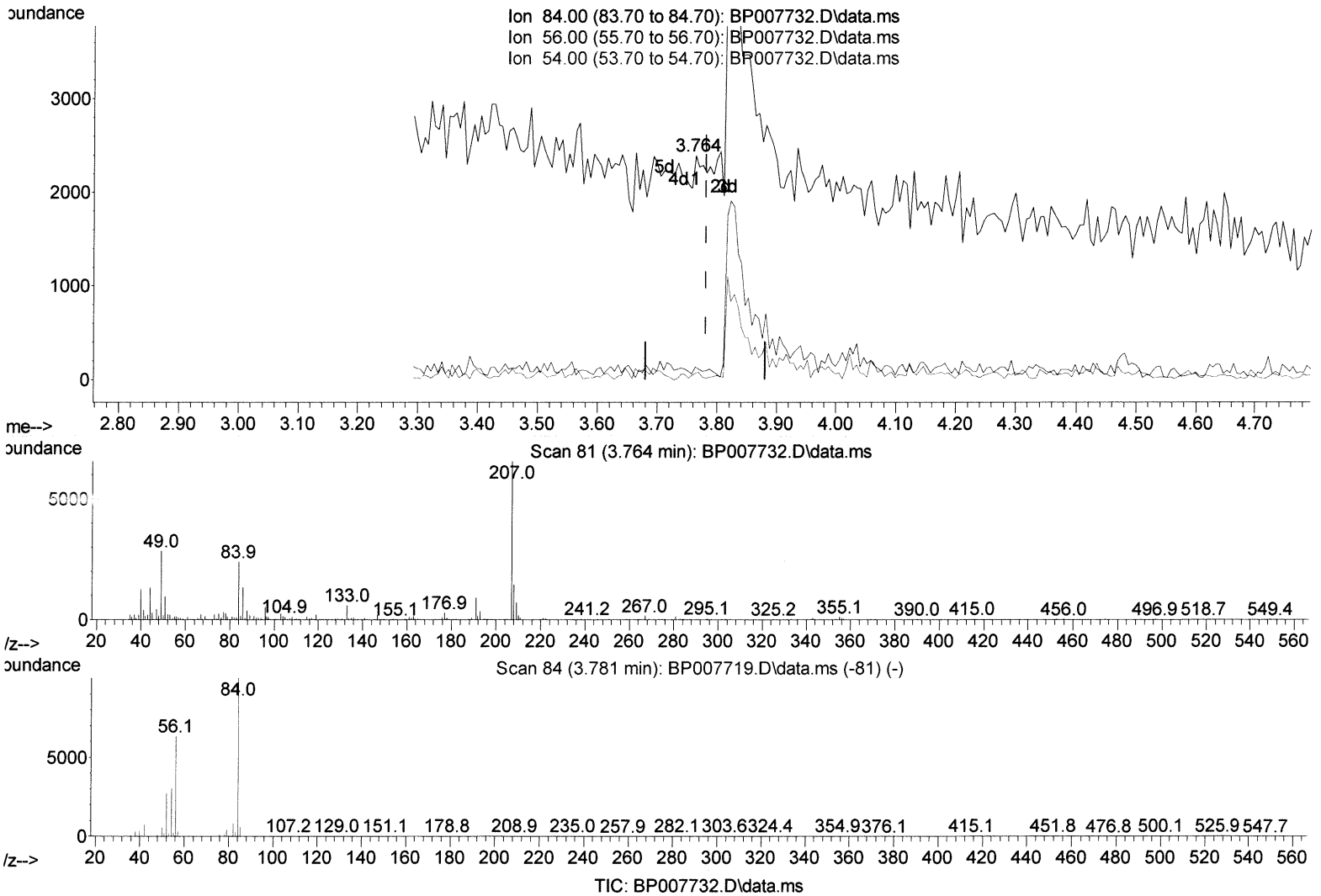


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(4) Pyridine-d5 (S)

3.764min (-0.017) 0.03 ng/u1

response	491	
Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	5.64#
54.00	31.70	2.88#
0.00	0.00	0.00

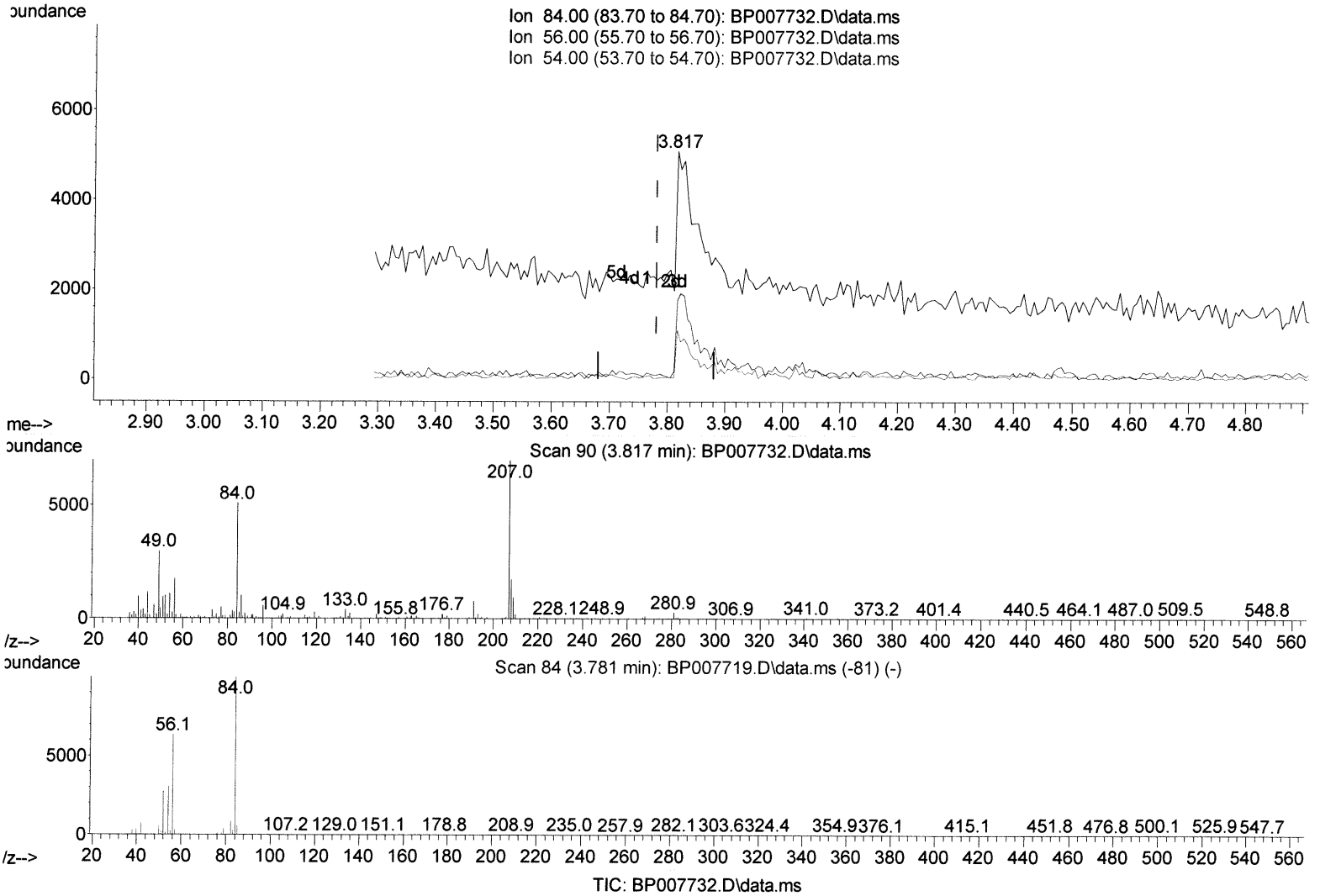
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(4) Pyridine-d5 (S)

3.817min (+ 0.036) 0.49 ng/ul m 10/29/21 ju

response 8045

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	66.30	34.52#
54.00	31.70	21.60#
0.00	0.00	0.00

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Manual Integrations
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	246530	20.000	ng/ul	0.00
20) Naphthalene-d8	10.734	136	993587	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	641179	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	1374283	20.000	ng/ul	0.00
79) Chrysene-d12	21.416	240	1488221	20.000	ng/ul	0.00
88) Perylene-d12	23.857	264	1577377	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	26117	4.812	ng/uL	0.00
4) Pyridine-d5	3.817	84	8045m	0.490	ng/ul	> 0.04 10/28/21 JU
7) Phenol-d5	7.093	99	639476	34.151	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	376868	36.250	ng/ul	0.00
11) 2-Chlorophenol-d4	7.464	132	514263	35.311	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	503238	34.321	ng/ul	0.00
21) Nitrobenzene-d5	9.087	128	241680	35.693	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	255190	34.507	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.352	165	493158	32.762	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	612378	30.767	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1533768	35.314	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	1841543	34.594	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	231084	29.026	ng/ul	0.00
60) Fluorene-d10	15.563	176	1332305	36.387	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	184605	25.113	ng/ul	0.00
73) Anthracene-d10	17.422	188	2084240	36.904	ng/ul	0.00
81) Pyrene-d10	19.657	212	2540044	36.660	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	2574593	33.226	ng/ul	-0.01

Target Compounds Qvalue

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