

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082024\
 Data File : BM047277.D
 Acq On : 20 Aug 2024 19:35
 Operator : RC/JU
 Sample : P3637-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ETGI - 370

Quant Time: Aug 21 02:48:57 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Aug 11 23:47:58 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.357	152	270066	20.000	ng	-0.02
21) Naphthalene-d8	10.116	136	933360	20.000	ng	-0.02
39) Acenaphthene-d10	14.016	164	565512	20.000	ng	-0.02
64) Phenanthrene-d10	16.780	188	1071047	20.000	ng	-0.01
76) Chrysene-d12	21.021	240	884079	20.000	ng	-0.01
86) Perylene-d12	23.721	264	908458	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.016	112	1864942	123.390	ng	0.00
7) Phenol-d6	6.575	99	2173541	104.367	ng	-0.01
23) Nitrobenzene-d5	8.510	82	1692703	91.319	ng	-0.01
42) 2,4,6-Tribromophenol	15.527	330	926529	141.299	ng	-0.01
45) 2-Fluorobiphenyl	12.627	172	3610329	98.478	ng	-0.02
79) Terphenyl-d14	19.451	244	4467314	108.275	ng	-0.01

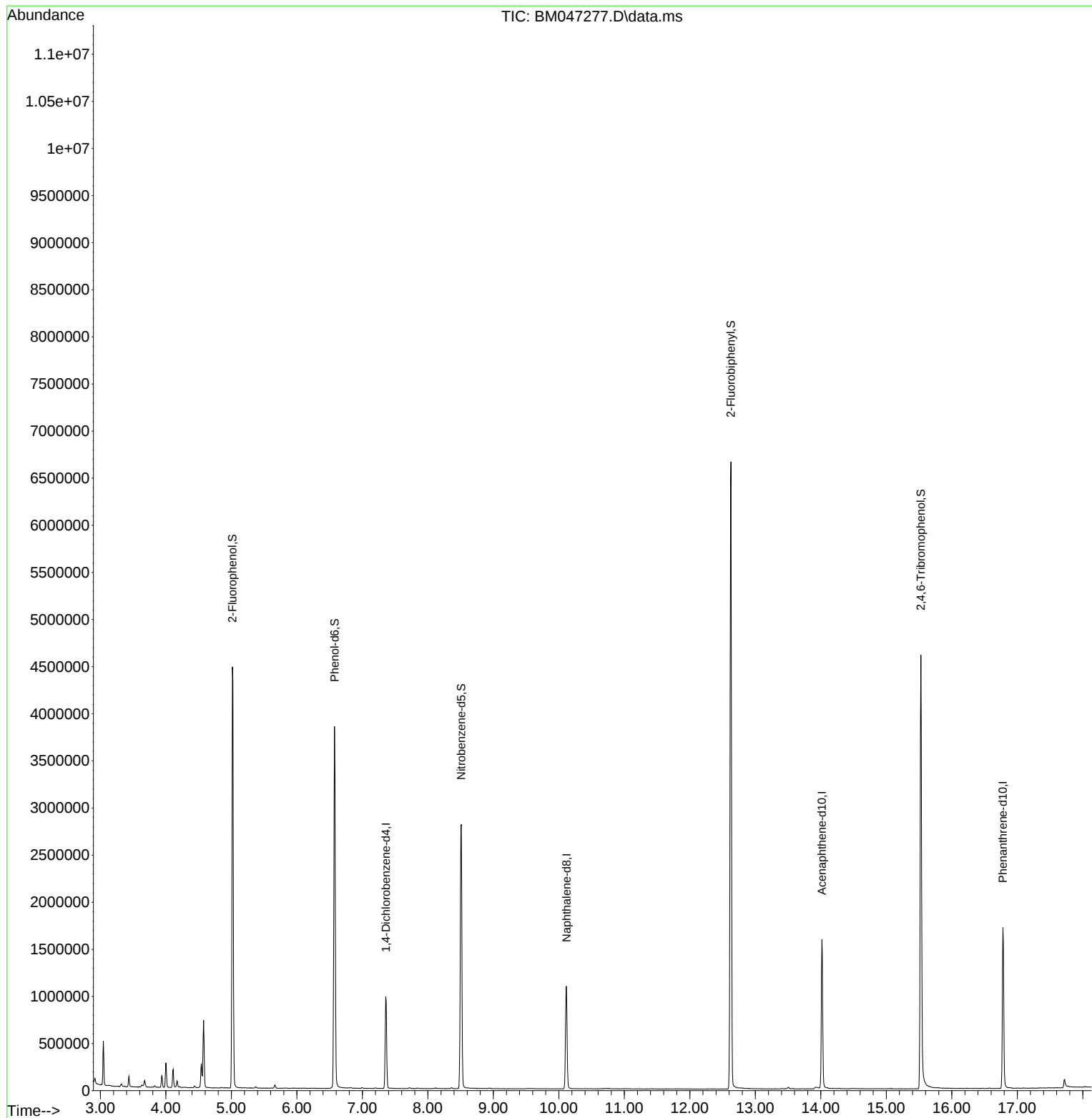
Target Compounds	Qvalue

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

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