

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082124\
 Data File : BM047295.D
 Acq On : 21 Aug 2024 13:07
 Operator : RC/JU
 Sample : P3617-08
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 VCPS-B4-081424-10-16

Quant Time: Aug 21 13:48:05 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	7.351	152	184312	20.000	ng	-0.02
21) Naphthalene-d8	10.110	136	689185	20.000	ng	-0.02
39) Acenaphthene-d10	14.016	164	448409	20.000	ng	-0.02
64) Phenanthrene-d10	16.780	188	1000271	20.000	ng	-0.01
76) Chrysene-d12	21.021	240	1043174	20.000	ng	-0.01
86) Perylene-d12	23.715	264	1051801	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.010	112	1123836	108.952	ng	-0.01
7) Phenol-d6	6.569	99	1474623	103.751	ng	-0.02
23) Nitrobenzene-d5	8.498	82	973494	71.125	ng	-0.02
42) 2,4,6-Tribromophenol	15.527	330	651936	125.387	ng	-0.01
45) 2-Fluorobiphenyl	12.622	172	2052614	70.610	ng	-0.02
79) Terphenyl-d14	19.445	244	3137401	64.445	ng	-0.02

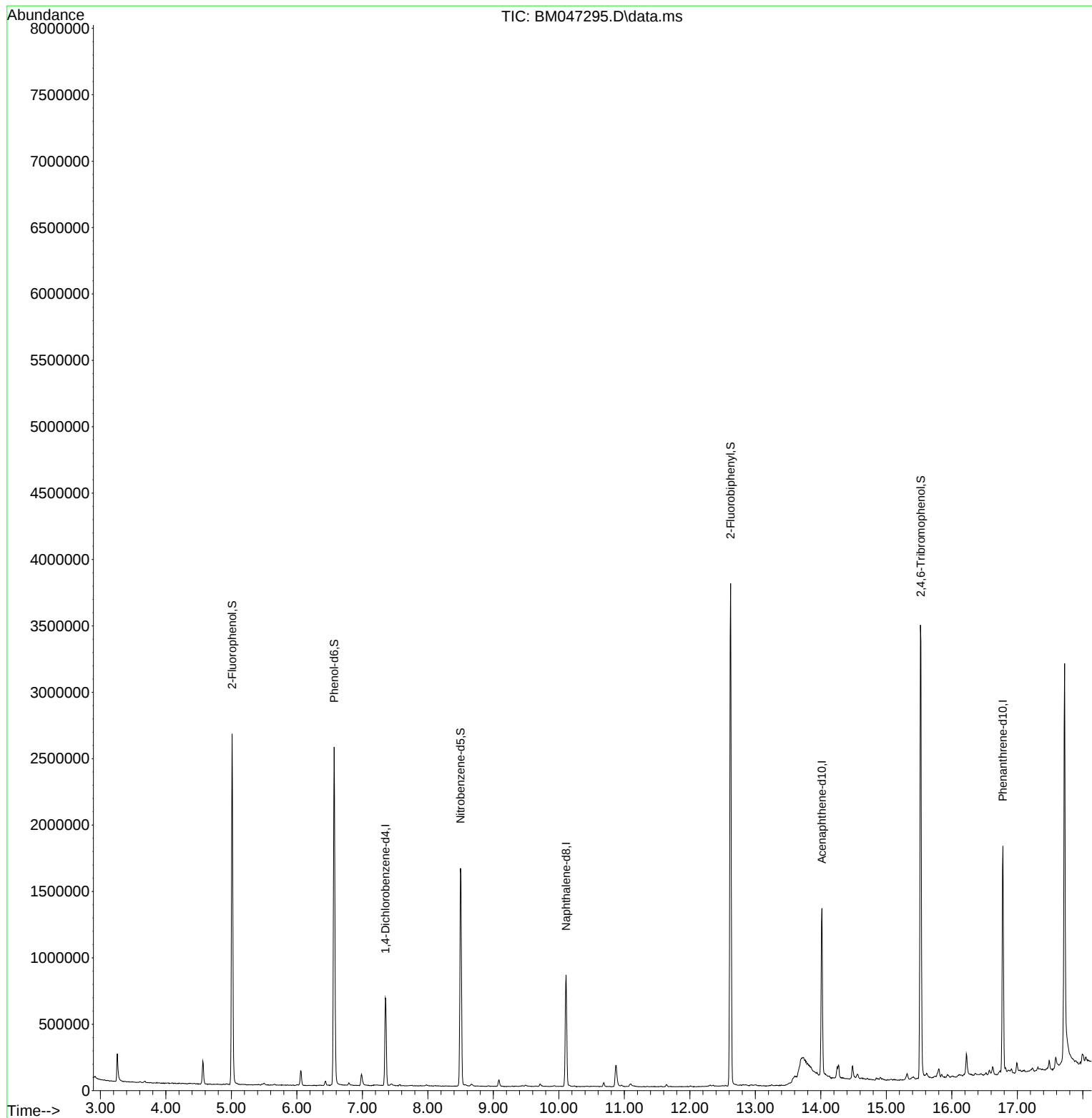
Target Compounds	Qvalue

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

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