

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110524\
Data File : BE101485.D
Acq On : 5 Nov 2024 18:39
Operator : RC/JU
Sample : P4643-12
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
TP-9

Quant Time: Nov 06 00:03:12 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE102824.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 29 01:26:30 2024
Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.567	152	60079	20.000	ng	0.00
21) Naphthalene-d8	10.334	136	257146	20.000	ng	0.00
39) Acenaphthene-d10	14.176	164	178286	20.000	ng	0.00
64) Phenanthrene-d10	16.914	188	428544	20.000	ng	0.00
76) Chrysene-d12	21.074	240	503216	20.000	ng	0.00
86) Perylene-d12	23.366	264	647563	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.322	112	478894	139.706	ng	0.00
7) Phenol-d6	6.944	99	573620	120.684	ng	0.00
23) Nitrobenzene-d5	8.777	82	387212	94.578	ng	0.00
42) 2,4,6-Tribromophenol	15.686	330	473043	151.384	ng	0.00
45) 2-Fluorobiphenyl	12.802	172	991641	94.196	ng	0.00
79) Terphenyl-d14	19.511	244	2259698	103.352	ng	0.00

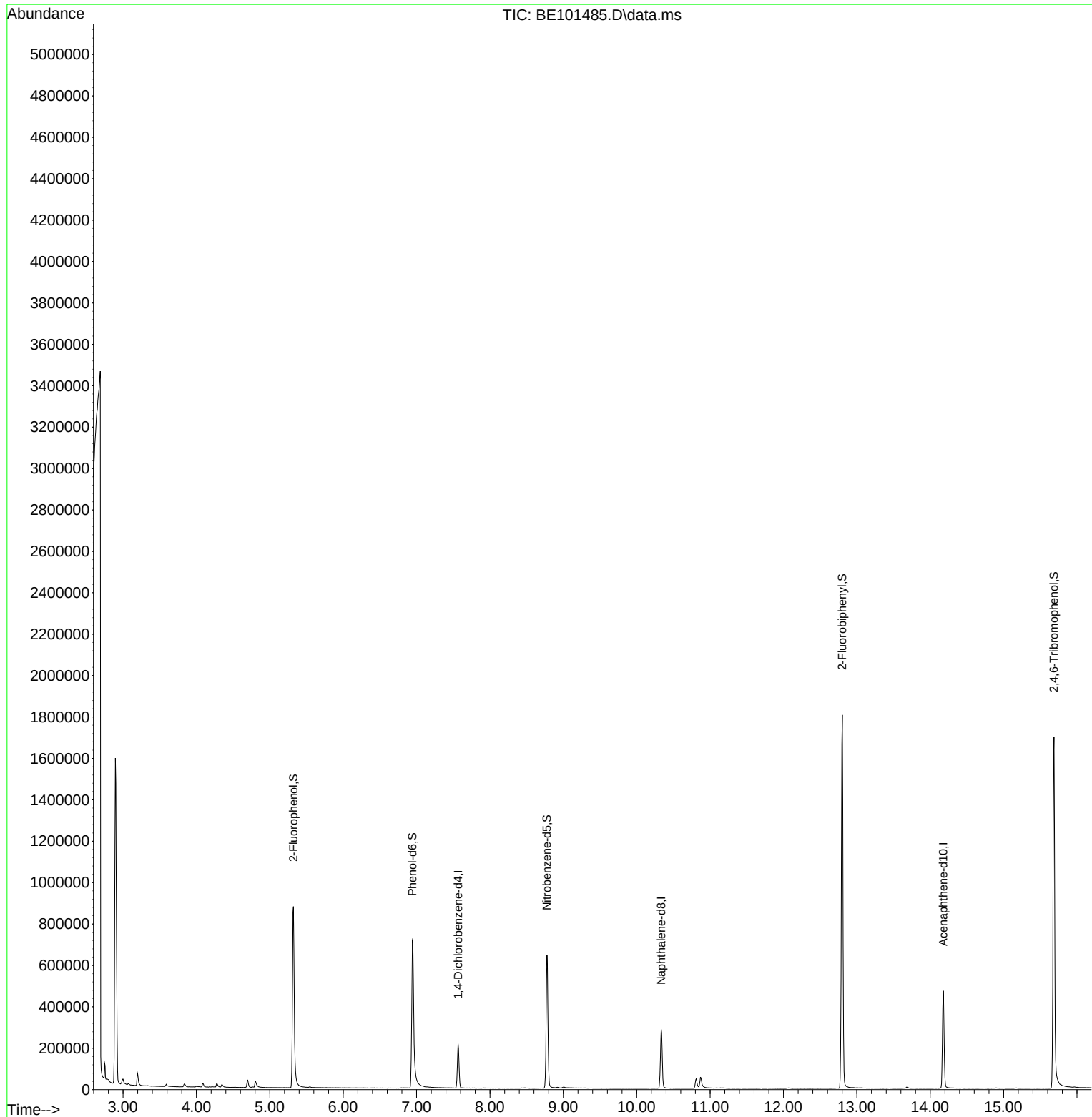
Target Compounds	Qvalue

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

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