

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE102924\
Data File : BE101413.D
Acq On : 29 Oct 2024 18:40
Operator : RC/JU
Sample : PB164369BL
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
PB164369BL

Quant Time: Oct 30 00:26:46 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	7.571	152	120768	20.000	ng	0.00
21) Naphthalene-d8	10.339	136	508333	20.000	ng	0.00
39) Acenaphthene-d10	14.175	164	326187	20.000	ng	0.00
64) Phenanthrene-d10	16.913	188	758793	20.000	ng	0.00
76) Chrysene-d12	21.073	240	995039	20.000	ng	0.00
86) Perylene-d12	23.359	264	1362650	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.321	112	1089923	158.177	ng	0.00
7) Phenol-d6	6.948	99	1356603	141.987	ng	0.00
23) Nitrobenzene-d5	8.782	82	781766	96.594	ng	0.00
42) 2,4,6-Tribromophenol	15.685	330	797601	139.514	ng	0.00
45) 2-Fluorobiphenyl	12.800	172	1979540	102.776	ng	0.00
79) Terphenyl-d14	19.510	244	3788985	87.640	ng	0.00

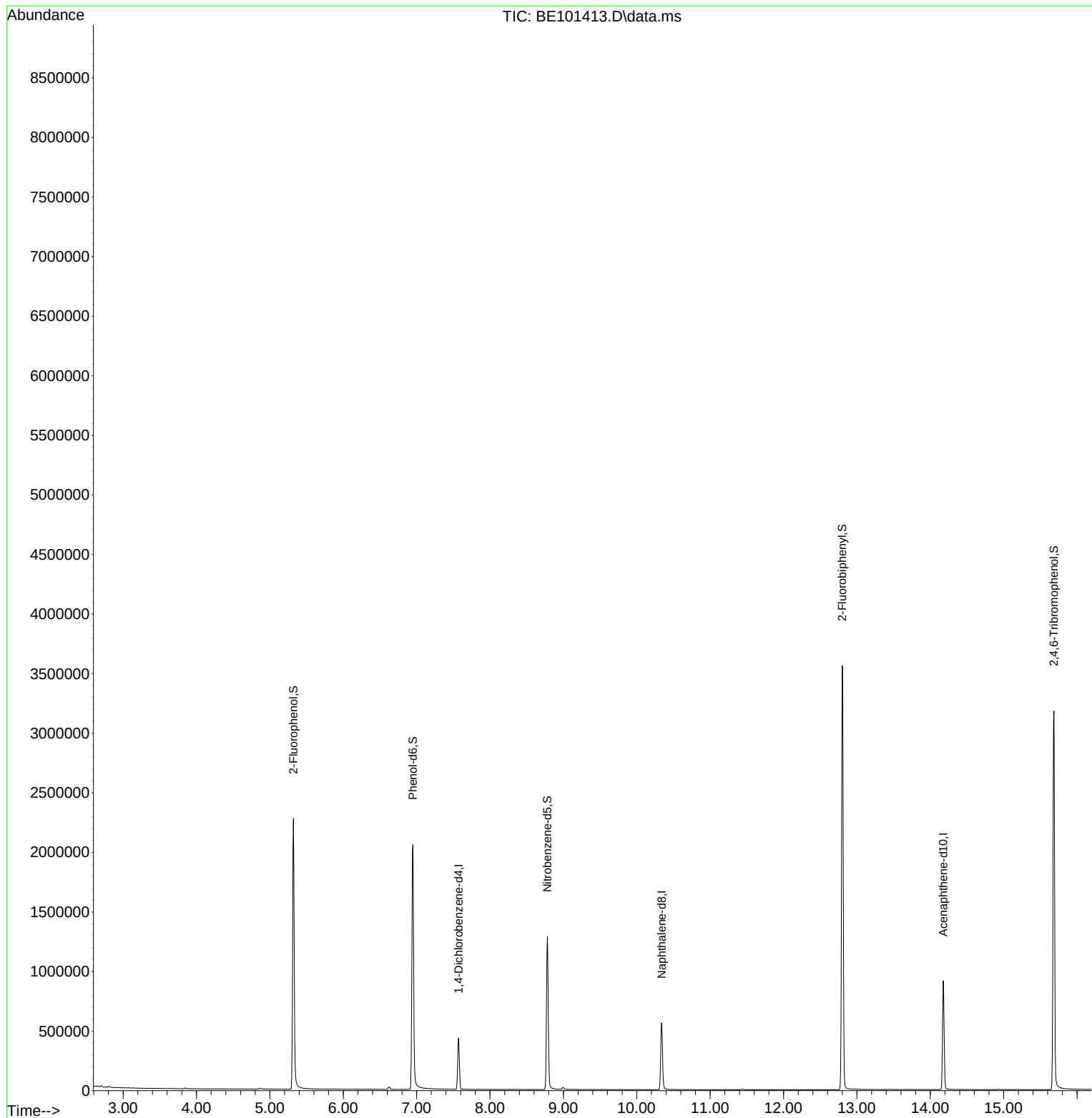
Target Compounds	Qvalue

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

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