

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE103024\
Data File : BE101422.D
Acq On : 30 Oct 2024 13:34
Operator : RC/JU
Sample : PB164382BL
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
PB164382BL

Quant Time: Oct 30 14:27:53 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	7.569	152	123851	20.000	ng	0.00
21) Naphthalene-d8	10.336	136	541935	20.000	ng	0.00
39) Acenaphthene-d10	14.179	164	358960	20.000	ng	0.00
64) Phenanthrene-d10	16.911	188	831647	20.000	ng	0.00
76) Chrysene-d12	21.071	240	1025552	20.000	ng	0.00
86) Perylene-d12	23.356	264	1388354	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.319	112	1128905	159.756	ng	0.00
7) Phenol-d6	6.946	99	1427387	145.676	ng	0.00
23) Nitrobenzene-d5	8.779	82	819469	94.975	ng	0.00
42) 2,4,6-Tribromophenol	15.683	330	883713	140.463	ng	0.00
45) 2-Fluorobiphenyl	12.804	172	2107705	99.439	ng	0.00
79) Terphenyl-d14	19.508	244	3901753	87.563	ng	0.00

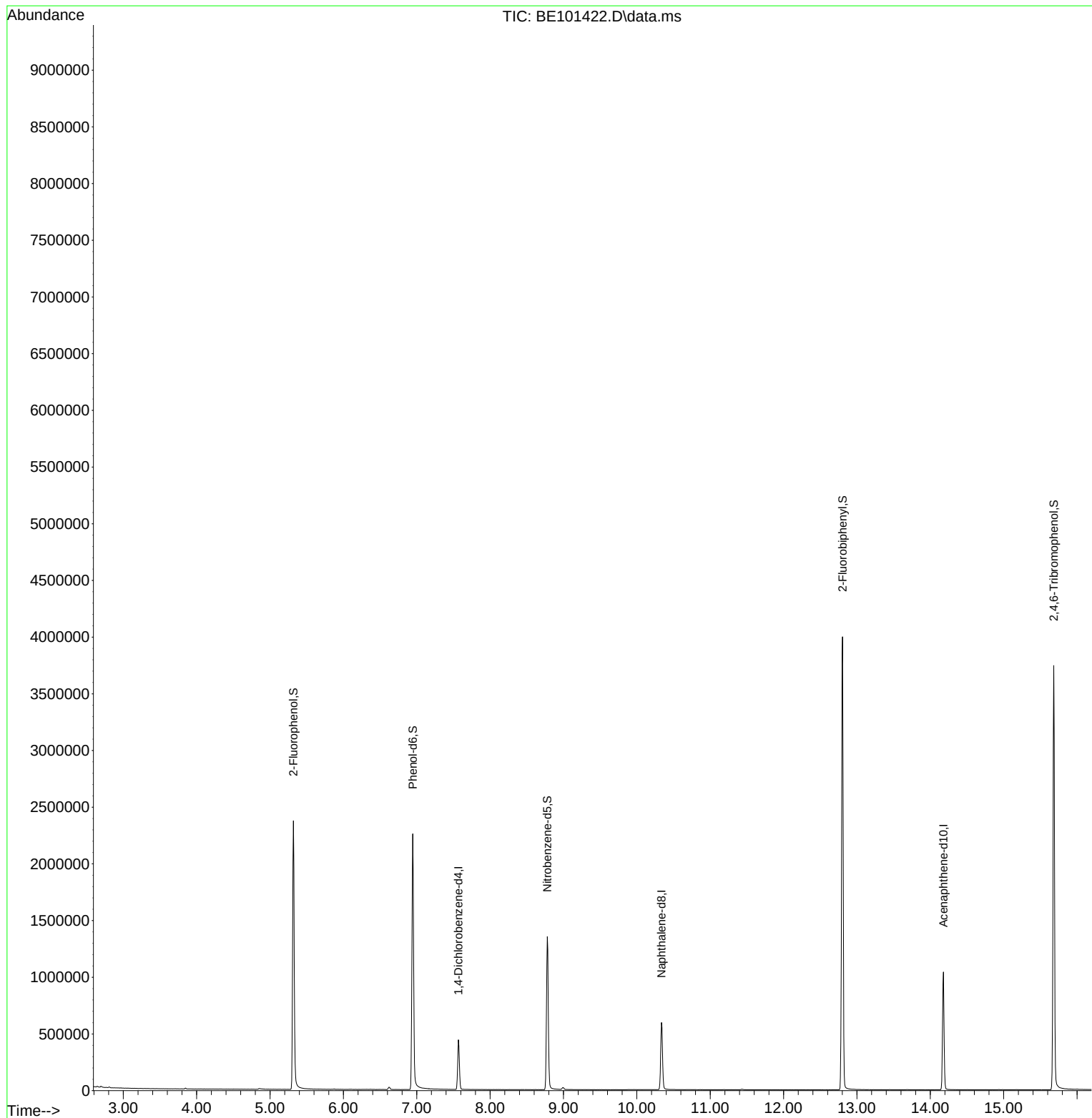
Target Compounds	Qvalue

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

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