

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110724\
 Data File : BE101527.D
 Acq On : 7 Nov 2024 12:38
 Operator : RC/JU
 Sample : PB164662TB
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB164662TB

Quant Time: Nov 07 12:28:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 07 00:01:20 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.564	152	53427	20.000	ng	0.00
21) Naphthalene-d8	10.331	136	209075	20.000	ng	0.00
39) Acenaphthene-d10	14.174	164	125183	20.000	ng	0.00
64) Phenanthrene-d10	16.912	188	279160	20.000	ng	0.00
76) Chrysene-d12	21.072	240	351169	20.000	ng	0.00
86) Perylene-d12	23.357	264	469581	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.319	112	520211	172.165	ng	0.00
7) Phenol-d6	6.947	99	625288	150.933	ng	0.00
23) Nitrobenzene-d5	8.774	82	377263	113.977	ng	0.00
42) 2,4,6-Tribromophenol	15.684	330	376554	159.857	ng	0.00
45) 2-Fluorobiphenyl	12.793	172	900559	117.463	ng	0.00
79) Terphenyl-d14	19.503	244	1758638	110.854	ng	0.00

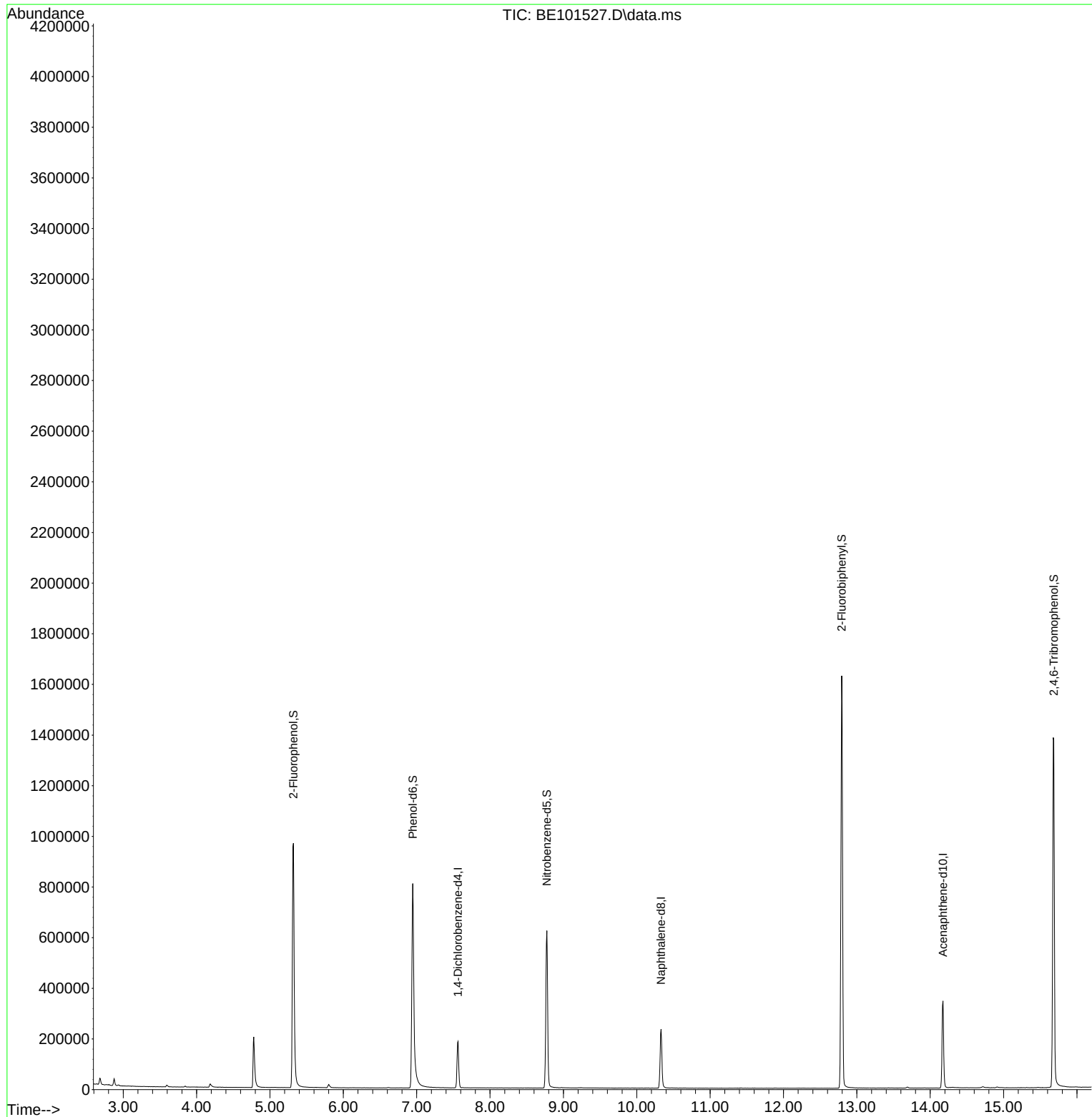
Target Compounds	Qvalue

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

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