

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE103124\
 Data File : BE101426.D
 Acq On : 31 Oct 2024 10:57
 Operator : RC/JU
 Sample : PB164533BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB164533BL

Quant Time: Oct 31 14:12:25 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	111412	20.000	ng	0.00
21) Naphthalene-d8	10.340	136	479387	20.000	ng	0.00
39) Acenaphthene-d10	14.177	164	322455	20.000	ng	0.00
64) Phenanthrene-d10	16.915	188	764639	20.000	ng	0.00
76) Chrysene-d12	21.075	240	978042	20.000	ng	0.00
86) Perylene-d12	23.361	264	1302952	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.317	112	1015593	159.767	ng	0.00
7) Phenol-d6	6.944	99	1307959	148.392	ng	0.00
23) Nitrobenzene-d5	8.783	82	750895	98.382	ng	0.00
42) 2,4,6-Tribromophenol	15.687	330	850212	150.437	ng	0.00
45) 2-Fluorobiphenyl	12.802	172	1974331	103.692	ng	0.00
79) Terphenyl-d14	19.512	244	3819837	89.889	ng	0.00

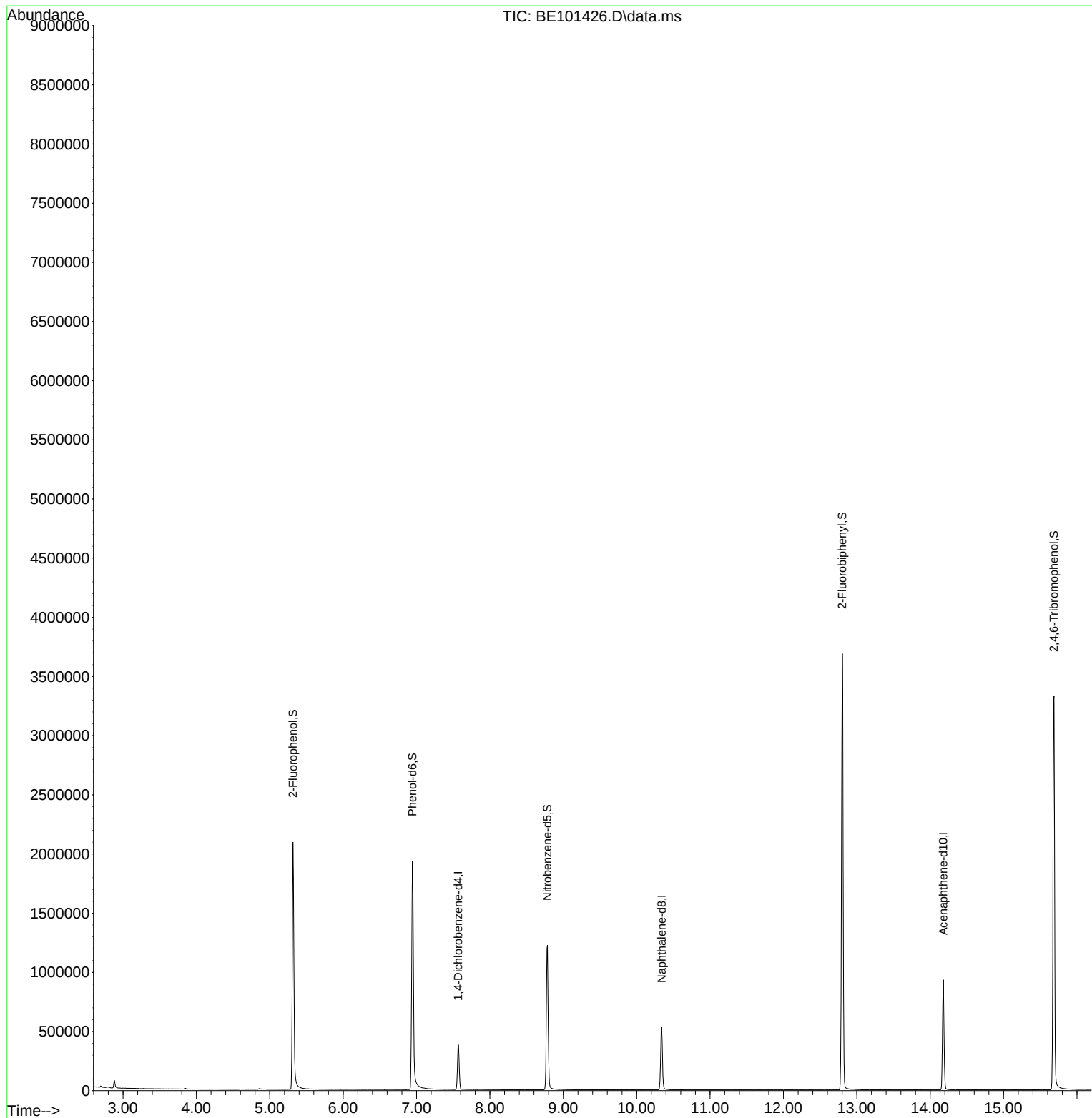
Target Compounds	Qvalue

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

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