

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110524\
 Data File : BE101486.D
 Acq On : 5 Nov 2024 19:15
 Operator : RC/JU
 Sample : P4682-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 BELL-SHOP-RAGS

Quant Time: Nov 06 00:03:29 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)

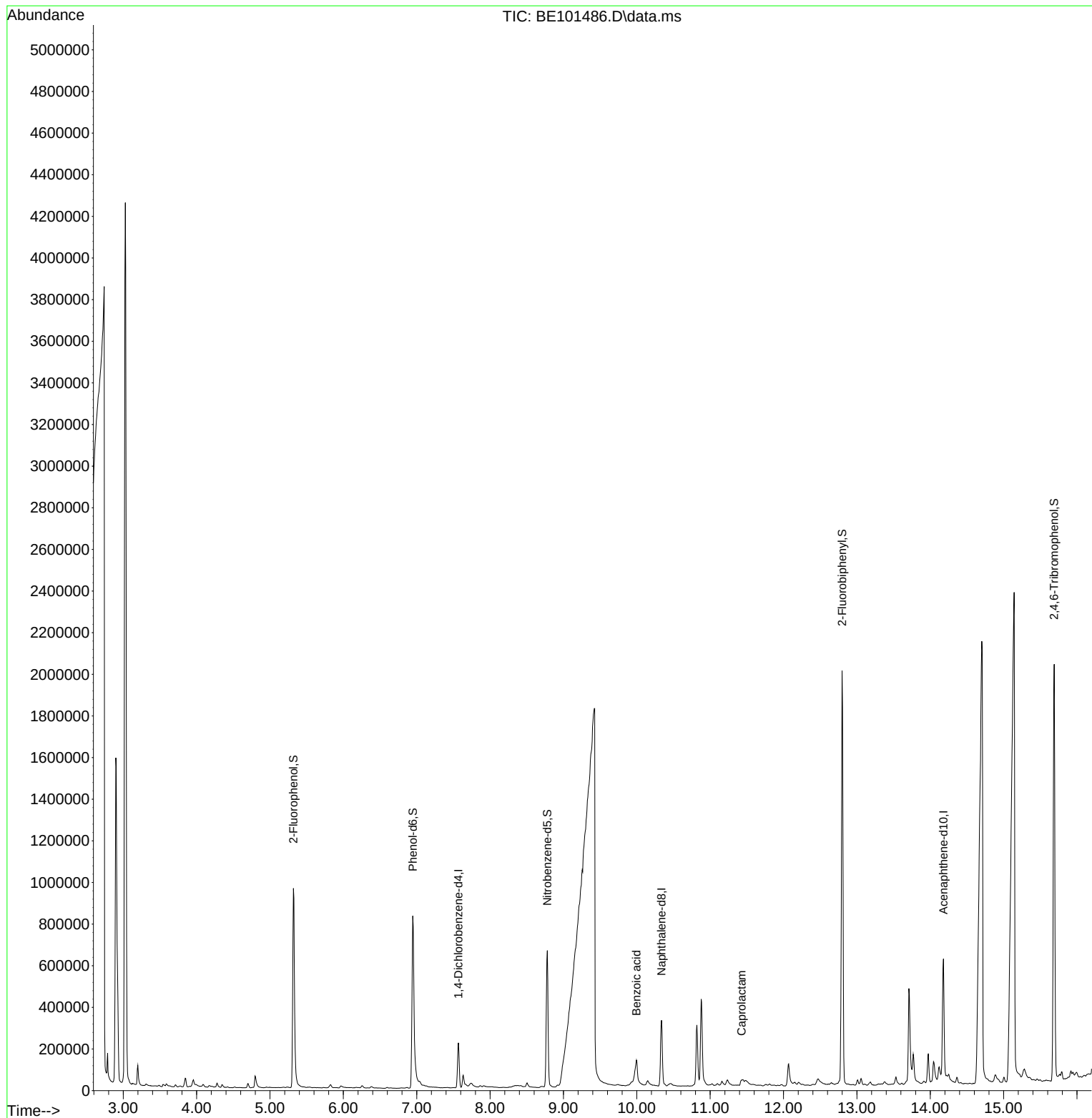
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.569	152	62298	20.000	ng	0.00
21) Naphthalene-d8	10.337	136	283913	20.000	ng	0.00
39) Acenaphthene-d10	14.179	164	202540	20.000	ng	0.00
64) Phenanthrene-d10	16.917	188	480864	20.000	ng	0.00
76) Chrysene-d12	21.077	240	520965	20.000	ng	0.00
86) Perylene-d12	23.374	264	659503	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.319	112	509097	143.227	ng	0.00
7) Phenol-d6	6.947	99	668526	135.641	ng	0.00
23) Nitrobenzene-d5	8.780	82	391402	86.589	ng	0.00
42) 2,4,6-Tribromophenol	15.689	330	520250	146.555	ng	0.00
45) 2-Fluorobiphenyl	12.799	172	1058488	88.505	ng	0.00
79) Terphenyl-d14	19.514	244	2158494	95.359	ng	0.00
Target Compounds						
						Qvalue
32) Benzoic acid	9.996	122	80347	30.340	ng	98
35) Caprolactam	11.430	113	6081	3.919	ng	91

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

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