

Data Path : Z:\svoasrv\HPCHEM1\BNA\_E\Data\BE111224\  
 Data File : BE101579.D  
 Acq On : 12 Nov 2024 13:05  
 Operator : RC/JU  
 Sample : P4722-10  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_E  
 ClientSampleId :  
 WC-2(0-6)

Quant Time: Nov 12 13:45:22 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

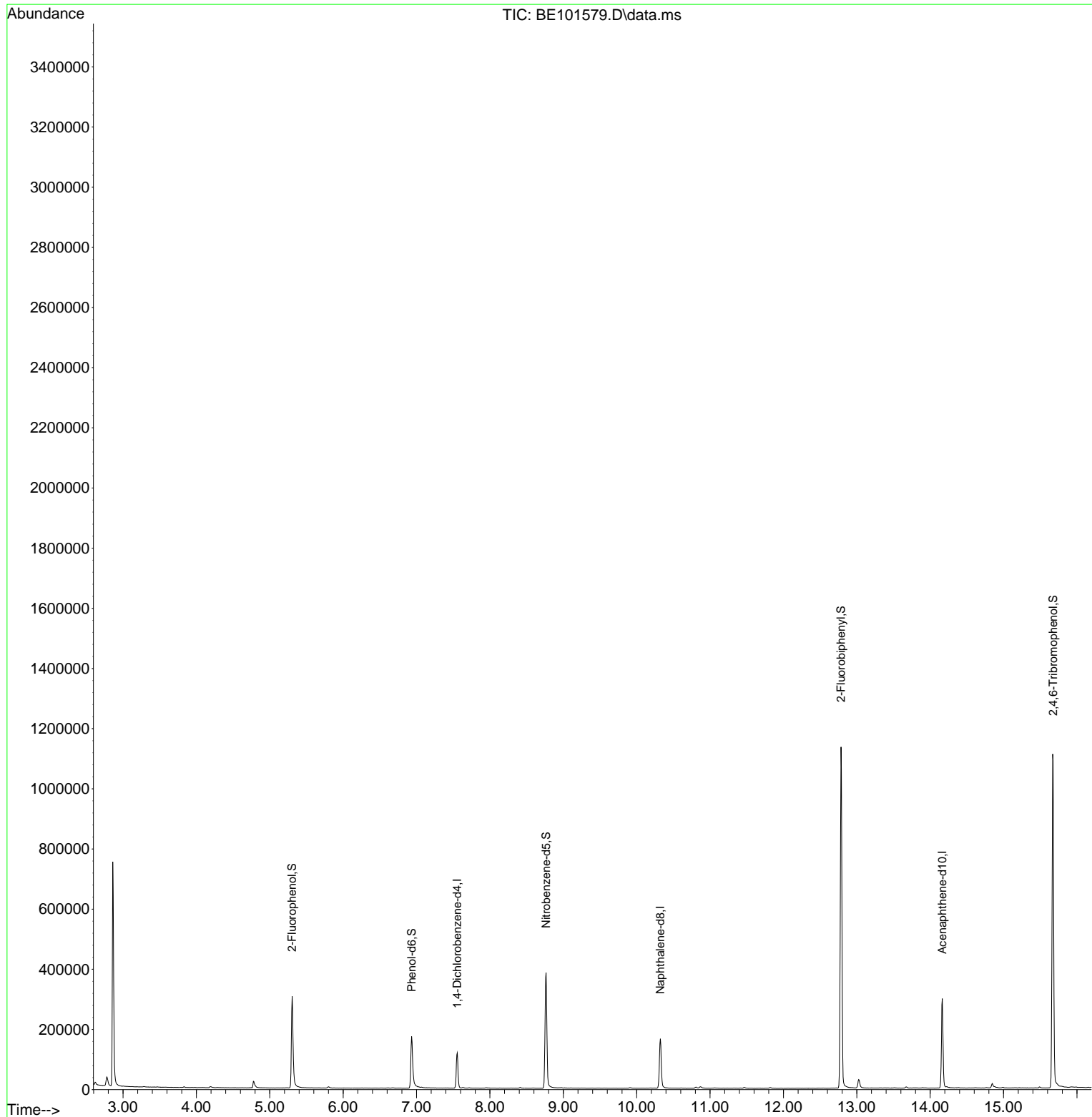
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.555	152	34307	20.000	ng	0.00
21) Naphthalene-d8	10.323	136	153776	20.000	ng	0.00
39) Acenaphthene-d10	14.165	164	106805	20.000	ng	0.00
64) Phenanthrene-d10	16.903	188	255892	20.000	ng	0.00
76) Chrysene-d12	21.063	240	316321	20.000	ng	-0.01
86) Perylene-d12	23.343	264	410447	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.305	112	156074	80.440	ng	0.00
7) Phenol-d6	6.932	99	128123	48.162	ng	-0.01
23) Nitrobenzene-d5	8.766	82	241488	99.193	ng	0.00
42) 2,4,6-Tribromophenol	15.675	330	285948	142.281	ng	0.00
45) 2-Fluorobiphenyl	12.784	172	625754	95.663	ng	-0.01
79) Terphenyl-d14	19.500	244	1562752	109.359	ng	0.00
Target Compounds						Qvalue
70) Pentachlorophenol	16.603	266	19106	9.480	ng	92

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

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