

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM102524\  
 Data File : BM048229.D  
 Acq On : 25 Oct 2024 09:28  
 Operator : RC/JU  
 Sample : PB164382BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

**Instrument :**  
 BNA\_M  
**ClientSampleId :**  
 PB164382BL

Quant Time: Oct 25 11:01:19 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM102324.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Oct 23 18:21:59 2024  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	221062	20.000	ng	0.00
21) Naphthalene-d8	10.510	136	804099	20.000	ng	0.00
39) Acenaphthene-d10	14.363	164	478139	20.000	ng	0.00
64) Phenanthrene-d10	17.104	188	895440	20.000	ng	0.00
76) Chrysene-d12	21.333	240	702826	20.000	ng	0.00
86) Perylene-d12	24.280	264	737871	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.304	112	1946989	148.048	ng	0.00
7) Phenol-d6	6.898	99	2344528	136.891	ng	0.00
23) Nitrobenzene-d5	8.875	82	1359612	95.470	ng	0.00
42) 2,4,6-Tribromophenol	15.851	330	759621	129.795	ng	0.00
45) 2-Fluorobiphenyl	12.980	172	2996789	96.176	ng	0.00
79) Terphenyl-d14	19.733	244	3663504	106.374	ng	0.00

Target Compounds Qvalue

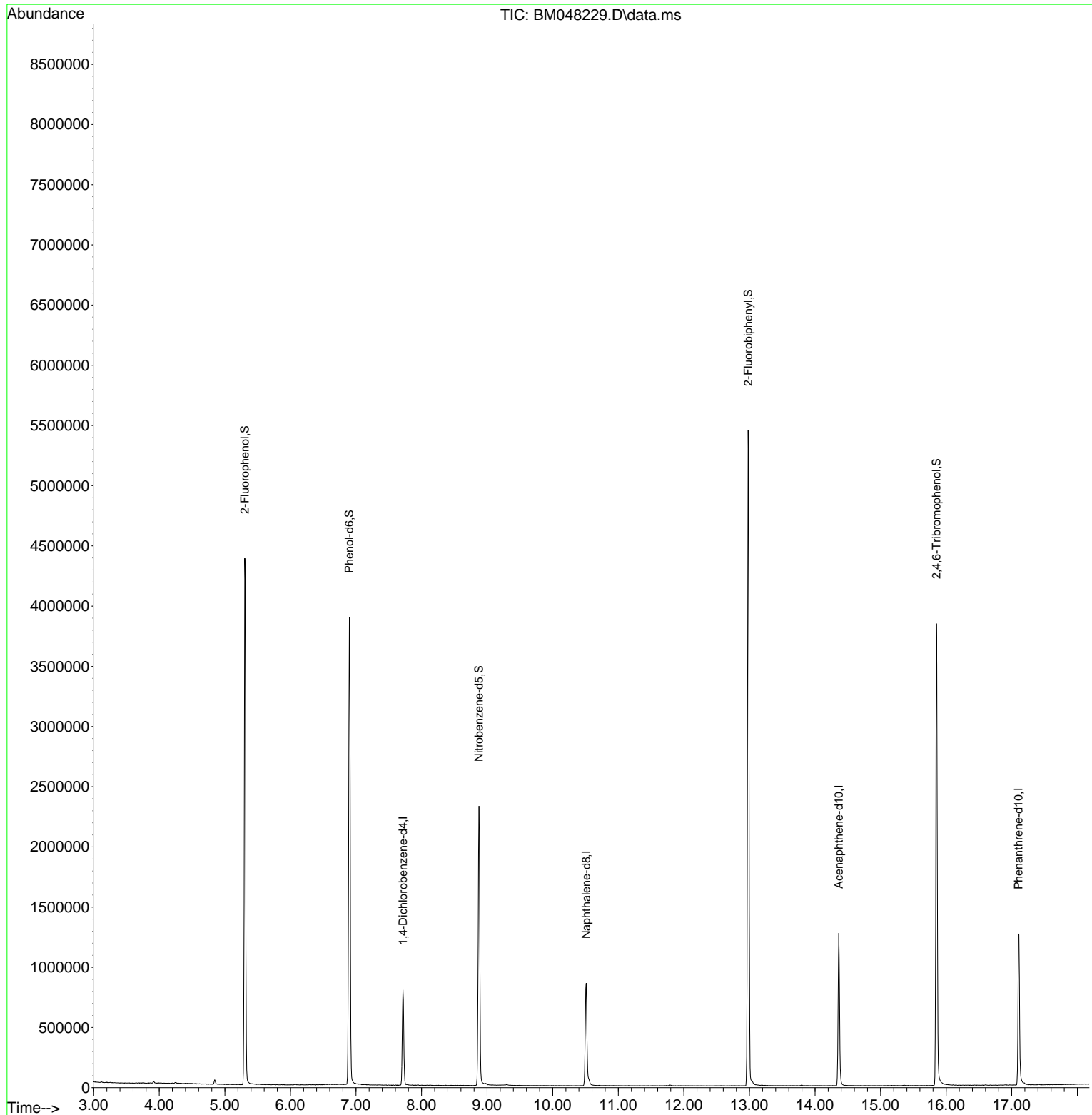
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

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