

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081924\
 Data File : BM047260.D
 Acq On : 19 Aug 2024 20:09
 Operator : RC/JU
 Sample : P3637-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ETGI - 370

Quant Time: Aug 20 02:01:59 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Aug 11 23:47:58 2024
 Response via : Initial Calibration

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

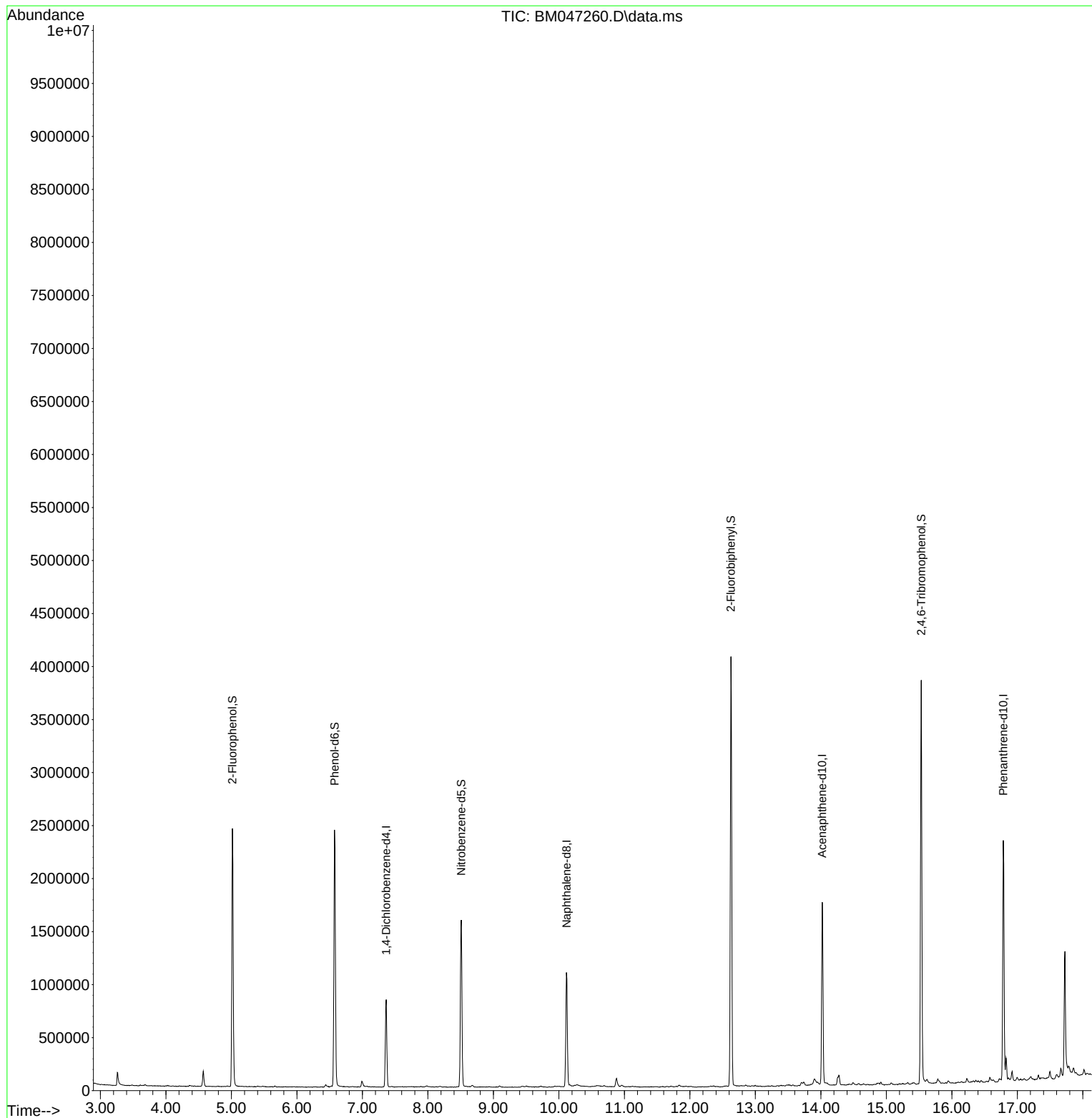
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.363	152	227341	20.000	ng	-0.01
21) Naphthalene-d8	10.116	136	888897	20.000	ng	-0.02
39) Acenaphthene-d10	14.021	164	608006	20.000	ng	-0.01
64) Phenanthrene-d10	16.786	188	1348748	20.000	ng	0.00
76) Chrysene-d12	21.027	240	1294513	20.000	ng	0.00
86) Perylene-d12	23.732	264	1242528	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.016	112	1056710	83.055	ng	0.00
7) Phenol-d6	6.575	99	1421425	81.079	ng	-0.01
23) Nitrobenzene-d5	8.510	82	918297	52.019	ng	-0.01
42) 2,4,6-Tribromophenol	15.533	330	707492	100.354	ng	0.00
45) 2-Fluorobiphenyl	12.627	172	2158113	54.752	ng	-0.02
79) Terphenyl-d14	19.456	244	3803196	62.953	ng	0.00
Target Compounds						
						Qvalue
75) Fluoranthene	18.862	202	400840	4.869	ng	98
78) Pyrene	19.233	202	370642	4.010	ng	99
81) Benzo(a)anthracene	21.009	228	211265	2.458	ng	99
83) Chrysene	21.068	228	221132	2.712	ng	98
84) Bis(2-ethylhexyl)phtha...	20.968	149	1144111	21.387	ng	100
88) Benzo(b)fluoranthene	22.885	252	273274	3.703	ng	# 97
90) Benzo(a)pyrene	23.603	252	163970	2.565	ng	# 90

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

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