

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040225\
 Data File : BM049787.D
 Acq On : 02 Apr 2025 12:49
 Operator : RC/JU
 Sample : Q1679-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-10

Quant Time: Apr 02 13:33:44 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031325.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 13 14:55:34 2025
 Response via : Initial Calibration

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

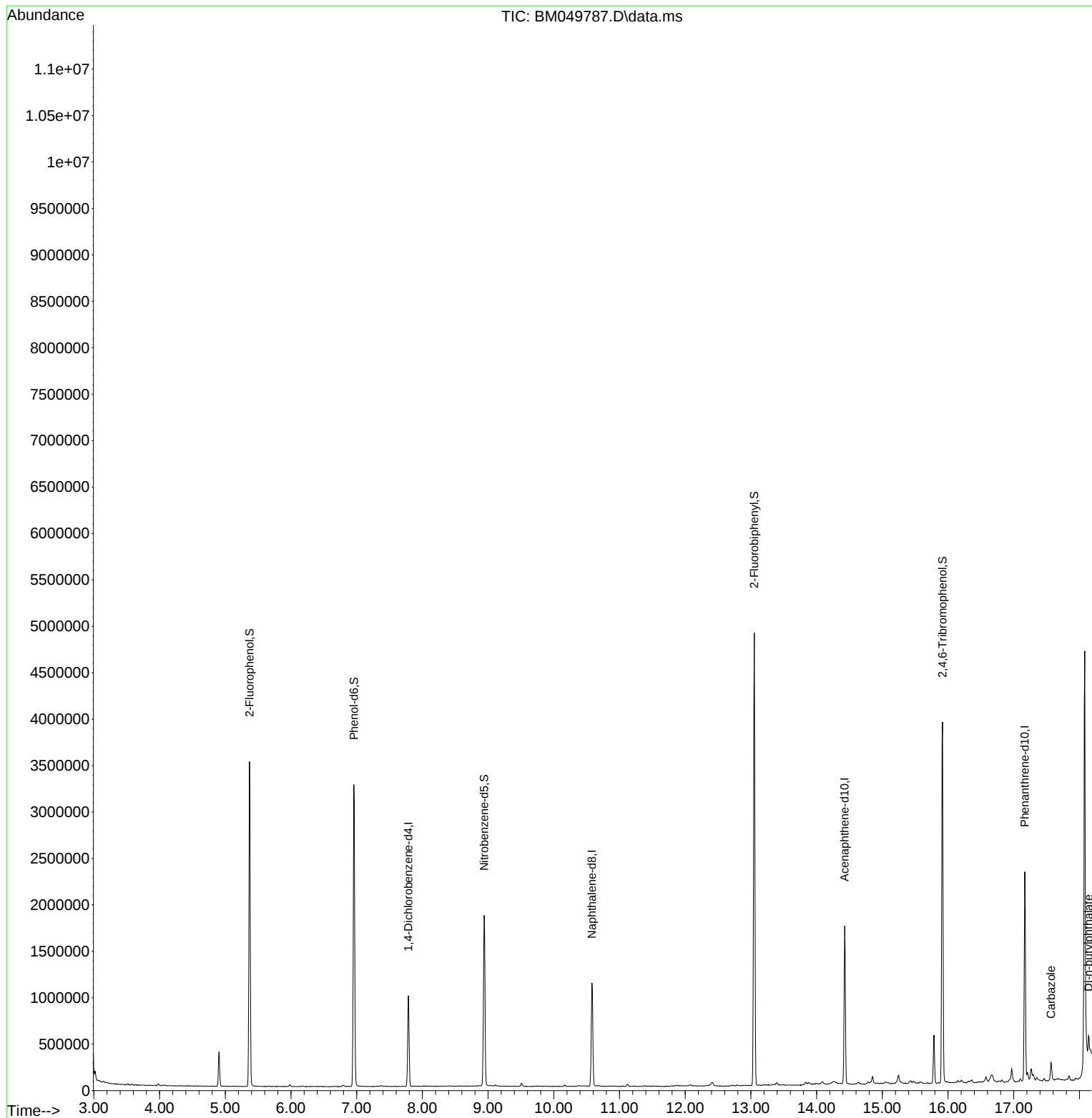
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	260998	20.000	ng	-0.02
21) Naphthalene-d8	10.580	136	891000	20.000	ng	-0.02
39) Acenaphthene-d10	14.427	164	573758	20.000	ng	-0.02
64) Phenanthrene-d10	17.168	188	1252828	20.000	ng	-0.02
76) Chrysene-d12	21.403	240	1706504	20.000	ng	-0.02
86) Perylene-d12	24.397	264	1701596	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	1369023	88.506	ng	-0.02
7) Phenol-d6	6.957	99	1698777	87.513	ng	-0.02
23) Nitrobenzene-d5	8.939	82	1025913	59.116	ng	-0.02
42) 2,4,6-Tribromophenol	15.915	330	669749	99.966	ng	-0.02
45) 2-Fluorobiphenyl	13.051	172	2354930	51.403	ng	-0.02
79) Terphenyl-d14	19.792	244	3900016	38.511	ng	-0.02
Target Compounds						
						Qvalue
73) Carbazole	17.568	167	131640	2.158	ng	99
74) Di-n-butylphthalate	18.139	149	176919	2.646	ng	# 97
75) Fluoranthene	19.227	202	204270	2.550	ng	99
80) Butylbenzylphthalate	20.486	149	106450	2.877	ng	92
81) Benzo(a)anthracene	21.386	228	350061	3.071	ng	99
83) Chrysene	21.444	228	355213	3.288	ng	100
84) Bis(2-ethylhexyl)phtha...	21.315	149	175807	3.049	ng	97
85) Di-n-octyl phthalate	22.450	149	201474	2.248	ng	# 80
88) Benzo(b)fluoranthene	23.450	252	314609	2.745	ng	# 98
89) Benzo(k)fluoranthene	23.509	252	328561	2.913	ng	99
90) Benzo(a)pyrene	24.256	252	208612	2.167	ng	98

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

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