

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082324\
 Data File : BM047331.D
 Acq On : 23 Aug 2024 13:29
 Operator : RC/JU
 Sample : P3518-21RE
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OU4-TS-11-080724RE

Quant Time: Aug 23 14:26:08 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)

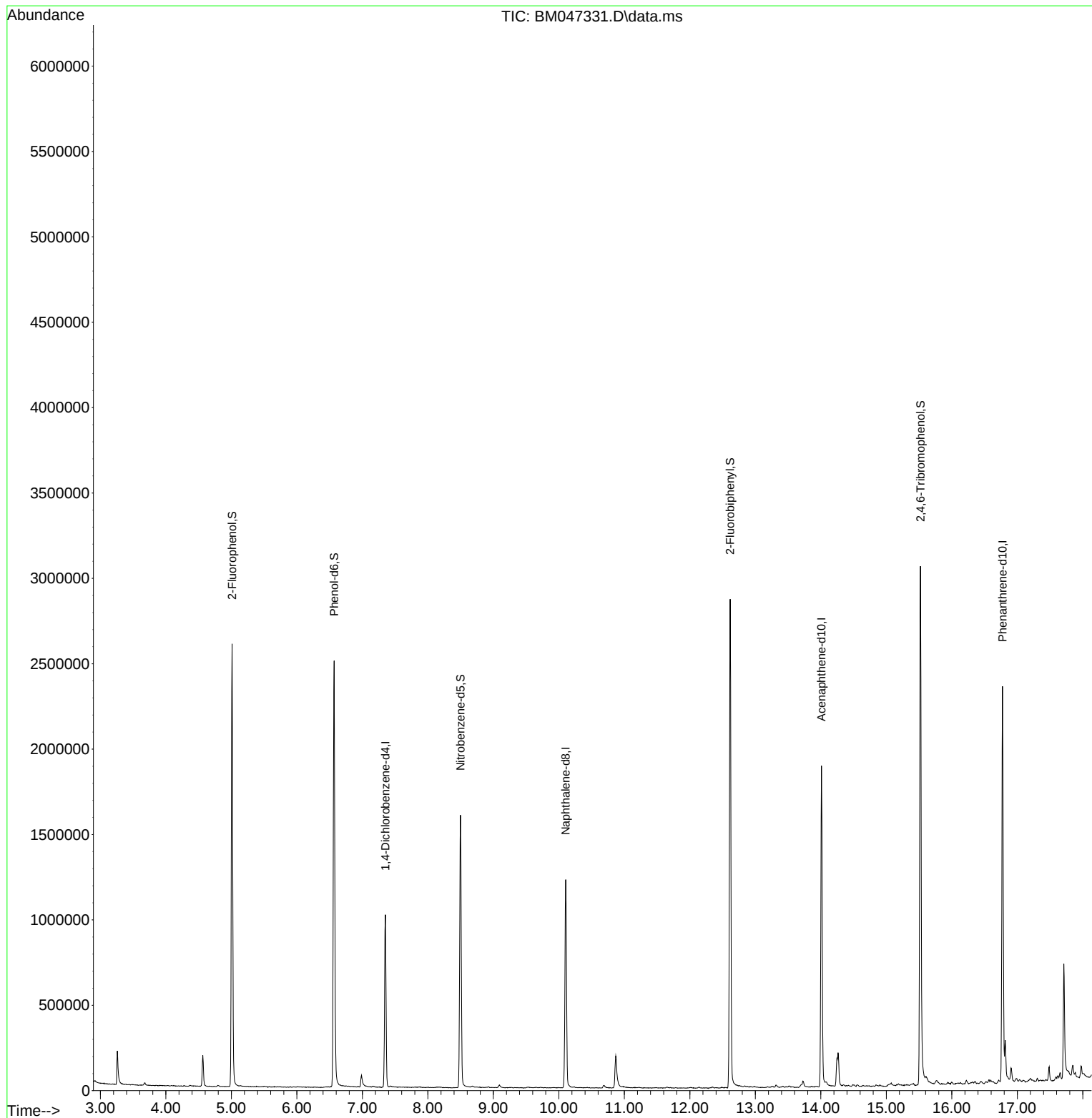
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.351	152	277753	20.000	ng	-0.02
21) Naphthalene-d8	10.104	136	1028963	20.000	ng	-0.03
39) Acenaphthene-d10	14.010	164	685603	20.000	ng	-0.02
64) Phenanthrene-d10	16.774	188	1451576	20.000	ng	-0.02
76) Chrysene-d12	21.015	240	1307088	20.000	ng	-0.02
86) Perylene-d12	23.703	264	1239770	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.010	112	1130035	72.697	ng	-0.01
7) Phenol-d6	6.569	99	1499381	70.003	ng	-0.02
23) Nitrobenzene-d5	8.498	82	961128	47.034	ng	-0.02
42) 2,4,6-Tribromophenol	15.521	330	623704	78.457	ng	-0.02
45) 2-Fluorobiphenyl	12.616	172	1619025	36.426	ng	-0.03
79) Terphenyl-d14	19.439	244	2477405	40.613	ng	-0.02
Target Compounds						
						Qvalue
75) Fluoranthene	18.851	202	459631	5.187	ng	99
78) Pyrene	19.221	202	405043	4.340	ng	99
81) Benzo(a)anthracene	20.998	228	200997	2.316	ng	97
83) Chrysene	21.051	228	214040	2.600	ng	98
88) Benzo(b)fluoranthene	22.862	252	244425	3.320	ng	# 95
90) Benzo(a)pyrene	23.580	252	149517	2.344	ng	97

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

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