

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060923\
 Data File : BM040214.D
 Acq On : 10 Jun 2023 06:42
 Operator : MA/JU
 Sample : 03046-07RE
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COMP-2-6-10RE

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 06/12/2023
 Supervised By :Jagrut Upadhyay 06/12/2023

Quant Time: Jun 10 08:35:54 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 08 23:25:50 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.987	152	515662	20.000	ng	0.00
21) Naphthalene-d8	10.798	136	2228301	20.000	ng	0.00
39) Acenaphthene-d10	14.627	164	1380203	20.000	ng	0.00
64) Phenanthrene-d10	17.368	188	1742648	20.000	ng	# 0.00
76) Chrysene-d12	21.592	240	394524m	20.000	ng	0.04
86) Perylene-d12	24.044	264	2183937	20.000	ng	0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.534	112	1916619	55.498	ng	0.00
7) Phenol-d6	7.140	99	1619169	34.360	ng	0.00
23) Nitrobenzene-d5	9.157	82	3089227	75.038	ng	0.00
42) 2,4,6-Tribromophenol	16.115	330	2295698	110.551	ng	0.00
45) 2-Fluorobiphenyl	13.251	172	8723003	83.533	ng	0.00
79) Terphenyl-d14	20.015	244	3364981	157.228	ng	0.03
Target Compounds						Qvalue
3) Pyridine	3.816	79	408817	10.393	ng	97
9) Benzaldehyde	7.116	77	75366	3.003	ng	98
22) Acetophenone	8.810	105	134267	2.307	ng	# 97

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

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