

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060823\
 Data File : BM040176.D
 Acq On : 09 Jun 2023 02:57
 Operator : MA/JU
 Sample : 03046-07
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 06/09/2023
 Supervised By :mohammad ahmed 06/13/2023

Quant Time: Jun 09 03:30:08 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	427492	20.000	ng	0.00
21) Naphthalene-d8	10.804	136	1609427	20.000	ng	0.00
39) Acenaphthene-d10	14.627	164	912804	20.000	ng	0.00
64) Phenanthrene-d10	17.368	188	1414223	20.000	ng	# 0.00
76) Chrysene-d12	21.592	240	374596m	20.000	ng	0.04
86) Perylene-d12	24.045	264	2098703	20.000	ng	0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.540	112	1919524	67.046	ng	0.00
7) Phenol-d6	7.146	99	1440690	36.878	ng	0.00
23) Nitrobenzene-d5	9.157	82	2437569	81.976	ng	0.00
42) 2,4,6-Tribromophenol	16.116	330	1706239	124.238	ng	0.00
45) 2-Fluorobiphenyl	13.257	172	5955130	86.228	ng	0.00
79) Terphenyl-d14	20.009	244	3057196	150.446	ng	0.02
Target Compounds						Qvalue
3) Pyridine	3.816	79	440833	13.518	ng	98
9) Benzaldehyde	7.122	77	63477	3.051	ng	96
22) Acetophenone	8.816	105	102769	2.445	ng	# 96

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

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