

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080822\
 Data File : BM036164.D
 Acq On : 09 Aug 2022 03:22
 Operator : CG/JU
 Sample : N4056-01DL 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TOTEDL

Quant Time: Aug 09 04:18:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 08 23:36:56 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 08/09/2022
 Supervised By :Jagrut Upadhyay 08/09/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.845	152	303323	20.000	ng	0.00
21) Naphthalene-d8	10.651	136	1173535	20.000	ng	0.00
39) Acenaphthene-d10	14.492	164	668838	20.000	ng	0.00
64) Phenanthrene-d10	17.233	188	1298871	20.000	ng	0.00
76) Chrysene-d12	21.415	240	1236169	20.000	ng	0.00
86) Perylene-d12	23.774	264	1368279	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	112960	6.038	ng	0.01
7) Phenol-d6	7.110	99	120964m	5.081	ng	0.05
23) Nitrobenzene-d5	9.016	82	140850	6.718	ng	0.00
42) 2,4,6-Tribromophenol	15.986	330	69482	8.855	ng	0.00
45) 2-Fluorobiphenyl	13.116	172	291269	6.430	ng	0.00
79) Terphenyl-d14	19.856	244	410129	6.544	ng	0.00
Target Compounds						Qvalue
9) Benzaldehyde	6.987	77	77464	6.146	ng	91
10) Phenol	7.140	94	88436	3.689	ng	93
20) 3+4-Methylphenols	8.663	107	1167896	54.176	ng	90
84) Bis(2-ethylhexyl)phtha...	21.327	149	421338	8.441	ng	# 98

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

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