

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112322\
 Data File : BP012716.D
 Acq On : 23 Nov 2022 18:23
 Operator : CG/JU
 Sample : N5749-16
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-83.5

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 11/25/2022
 Supervised By :Jagrut Upadhyay 11/25/2022

Quant Time: Nov 24 02:32:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 23 00:38:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.034	152	351977	20.000	ng	0.00
21) Naphthalene-d8	10.852	136	1313071	20.000	ng	0.00
39) Acenaphthene-d10	14.663	164	727317	20.000	ng	0.00
64) Phenanthrene-d10	17.422	188	1444857	20.000	ng	0.00
76) Chrysene-d12	21.504	240	1428997	20.000	ng	0.00
86) Perylene-d12	24.027	264	1591278	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	2874962	176.955	ng	0.00
7) Phenol-d6	7.181	99	3368384	144.566	ng	0.00
23) Nitrobenzene-d5	9.199	82	1977371	83.986	ng	0.00
42) 2,4,6-Tribromophenol	16.151	330	1266190	134.274	ng	0.00
45) 2-Fluorobiphenyl	13.293	172	4771749	93.468	ng	0.00
79) Terphenyl-d14	19.969	244	7348616	109.297	ng	0.00
Target Compounds						
32) Benzoic acid	10.040	122	35543m	7.735	ng	Qvalue
74) Di-n-butylphthalate	18.369	149	7313	5.787	ng	# 97

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

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