

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112421\
 Data File : BM033254.D
 Acq On : 24 Nov 2021 19:24
 Operator : CG/JU
 Sample : M4812-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SETTLING-TANK-01

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 11/30/2021
 Supervised By :Sohil Jodhani 11/30/2021

Quant Time: Nov 26 04:10:03 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM112421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 24 15:16:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.937	152	133201	20.000	ng	0.00
21) Naphthalene-d8	10.736	136	514452	20.000	ng	# 0.00
39) Acenaphthene-d10	14.560	164	303193	20.000	ng	0.00
64) Phenanthrene-d10	17.295	188	609014	20.000	ng	# 0.00
76) Chrysene-d12	21.453	240	624755	20.000	ng	0.00
86) Perylene-d12	23.789	264	657577	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.519	112	1025491	127.137	ng	0.01
7) Phenol-d6	7.107	99	1243754	115.133	ng	0.01
23) Nitrobenzene-d5	9.101	82	1010041	91.144	ng	0.00
42) 2,4,6-Tribromophenol	16.048	330	468178	141.654	ng	0.00
45) 2-Fluorobiphenyl	13.183	172	1984989	92.332	ng	0.00
79) Terphenyl-d14	19.906	244	2754169	87.171	ng	0.00
Target Compounds						
32) Benzoic acid	9.984	122	5903m	6.410	ng	Qvalue

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

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