

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022522\
 Data File : BM034088.D
 Acq On : 25 Feb 2022 21:26
 Operator : CG/JU
 Sample : N1664-11
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-4

Quant Time: Feb 26 04:02:49 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 26 03:45:21 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 02/28/2022
 Supervised By :Jagrut Upadhyay 02/28/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.001	152	118152	20.000	ng	0.00
21) Naphthalene-d8	10.819	136	419560	20.000	ng	# 0.00
39) Acenaphthene-d10	14.636	164	272537	20.000	ng	0.00
64) Phenanthrene-d10	17.377	188	567805	20.000	ng	# 0.00
76) Chrysene-d12	21.547	240	544793	20.000	ng	# 0.00
86) Perylene-d12	23.994	264	551832	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.554	112	935840	141.195	ng	0.00
7) Phenol-d6	7.154	99	992095	113.988	ng	0.00
23) Nitrobenzene-d5	9.166	82	757792	97.470	ng	0.00
42) 2,4,6-Tribromophenol	16.118	330	589721	178.287	ng	0.00
45) 2-Fluorobiphenyl	13.271	172	2057419	100.577	ng	0.00
79) Terphenyl-d14	19.995	244	3384334	108.174	ng	0.00
Target Compounds						
32) Benzoic acid	10.030	122	11411m	2.735	ng	Qvalue

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

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