

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031522\
 Data File : BM034249.D
 Acq On : 15 Mar 2022 22:01
 Operator : CG/JU
 Sample : PB143311BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB143311BL

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/16/2022
 Supervised By : Jagrut Upadhyay 03/16/2022

Quant Time: Mar 16 03:09:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 16 02:38:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.948	152	144454	20.000	ng	0.00
21) Naphthalene-d8	10.760	136	534607	20.000	ng	0.00
39) Acenaphthene-d10	14.589	164	319899	20.000	ng	0.00
64) Phenanthrene-d10	17.324	188	563423	20.000	ng	0.00
76) Chrysene-d12	21.495	240	405233m	20.000	ng	0.00
86) Perylene-d12	23.912	264	407753	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.502	112	1284184	159.482	ng	0.00
7) Phenol-d6	7.107	99	1586734	137.631	ng	0.00
23) Nitrobenzene-d5	9.113	82	1065780	97.061	ng	0.00
42) 2,4,6-Tribromophenol	16.071	330	553279	128.316	ng	0.00
45) 2-Fluorobiphenyl	13.219	172	2246974	95.982	ng	0.00
79) Terphenyl-d14	19.942	244	2512089	107.717	ng	0.00

Target Compounds Qvalue

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

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