

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052922\
 Data File : BP010617.D
 Acq On : 29 May 2022 14:05
 Operator : CG/JU
 Sample : N2859-08
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PE-D3

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 05/31/2022
 Supervised By :Jagrut Upadhyay 05/31/2022

Quant Time: May 30 02:06:06 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 28 02:18:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	110262	20.000	ng	0.00
21) Naphthalene-d8	10.669	136	421975	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	257845	20.000	ng	0.00
64) Phenanthrene-d10	17.257	188	532237	20.000	ng	0.00
76) Chrysene-d12	21.339	240	498392	20.000	ng	0.00
86) Perylene-d12	23.751	264	503645	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.452	112	399850	66.156	ng	0.00
7) Phenol-d6	7.040	99	303855	35.412	ng	0.00
23) Nitrobenzene-d5	9.028	82	819156	91.779	ng	0.00
42) 2,4,6-Tribromophenol	15.998	330	426154	146.009	ng	0.00
45) 2-Fluorobiphenyl	13.128	172	1671478	92.002	ng	0.00
79) Terphenyl-d14	19.810	244	2484244	93.625	ng	0.00
Target Compounds						Qvalue
32) Benzoic acid	9.963	122	6239m	8.700	ng	

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

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