

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030922\
 Data File : BP009337.D
 Acq On : 09 Mar 2022 22:39
 Operator : CG/JU
 Sample : N1845-04 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SAMPLE-4

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 03/10/2022
 Supervised By :Jagrut Upadhyay 03/10/2022

Quant Time: Mar 10 04:12:33 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 05 01:55:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	407016	20.000	ng	0.00
21) Naphthalene-d8	10.610	136	1804245	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	1266911	20.000	ng	0.00
64) Phenanthrene-d10	17.216	188	2997889	20.000	ng	0.00
76) Chrysene-d12	21.327	240	3462179	20.000	ng	0.00
86) Perylene-d12	23.733	264	3379011	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.410	112	431268	15.467	ng	0.00
7) Phenol-d6	6.987	99	1207503	34.077	ng	0.00
23) Nitrobenzene-d5	8.969	82	777207	23.726	ng	0.00
42) 2,4,6-Tribromophenol	15.951	330	90034	5.885	ng	0.00
45) 2-Fluorobiphenyl	13.081	172	2056336	22.321	ng	0.00
79) Terphenyl-d14	19.792	244	4548064	24.790	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	17.257	178	438916	2.548	ng	98
75) Fluoranthene	19.239	202	919186	4.622	ng	98
78) Pyrene	19.586	202	854615	3.656	ng	99
81) Benzo(a)anthracene	21.310	228	474886	2.024	ng	93
88) Benzo(b)fluoranthene	22.998	252	563403m	2.469	ng	

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

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