

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032322\
 Data File : BP009521.D
 Acq On : 23 Mar 2022 23:41
 Operator : CG/JU
 Sample : N2031-12 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CMP-B2-031822-0-2

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/24/2022
 Supervised By : Jagrut Upadhyay 03/24/2022

Quant Time: Mar 24 01:23:17 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 24 01:19:11 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	423016	20.000	ng	0.00
21) Naphthalene-d8	10.581	136	1543397	20.000	ng	0.00
39) Acenaphthene-d10	14.434	164	851163	20.000	ng	0.00
64) Phenanthrene-d10	17.192	188	1769550	20.000	ng	0.00
76) Chrysene-d12	21.310	240	1869649	20.000	ng	0.00
86) Perylene-d12	23.721	264	2045183	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	681783	28.140	ng	0.00
7) Phenol-d6	6.975	99	815864	24.900	ng	0.00
23) Nitrobenzene-d5	8.946	82	504518	17.371	ng	0.00
42) 2,4,6-Tribromophenol	15.934	330	284411	20.249	ng	0.00
45) 2-Fluorobiphenyl	13.051	172	1122882	18.288	ng	0.00
79) Terphenyl-d14	19.769	244	1726353	16.573	ng	-0.01
Target Compounds						Qvalue
78) Pyrene	19.569	202	412829	3.351	ng	96
81) Benzo(a)anthracene	21.298	228	251794	2.050	ng	# 67
83) Chrysene	21.345	228	490689	4.219	ng	# 83
88) Benzo(b)fluoranthene	22.986	252	260002m	2.130	ng	
90) Benzo(a)pyrene	23.610	252	279571	2.676	ng	# 79

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

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