

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033122\
 Data File : BP009647.D
 Acq On : 31 Mar 2022 19:32
 Operator : CG/JU
 Sample : N2202-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MHE

Quant Time: Apr 01 04:11:22 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 31 04:36:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.740	152	375079	20.000	ng	0.00
21) Naphthalene-d8	10.534	136	1490216	20.000	ng	0.00
39) Acenaphthene-d10	14.393	164	901892	20.000	ng	0.00
64) Phenanthrene-d10	17.157	188	1787970	20.000	ng	0.00
76) Chrysene-d12	21.275	240	1648625	20.000	ng	0.00
86) Perylene-d12	23.663	264	1827309	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.346	112	2194909	102.169	ng	0.00
7) Phenol-d6	6.934	99	2810348	96.733	ng	0.00
23) Nitrobenzene-d5	8.899	82	1801073	64.226	ng	0.00
42) 2,4,6-Tribromophenol	15.892	330	1176409	79.045	ng	0.00
45) 2-Fluorobiphenyl	13.010	172	4038745	62.079	ng	0.00
79) Terphenyl-d14	19.739	244	5561264	60.547	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	6.958	94	73978	2.538	ng	# 61
71) Phenanthrene	17.198	178	241887	2.527	ng	99
75) Fluoranthene	19.181	202	262537	2.266	ng	99
78) Pyrene	19.539	202	225659	2.077	ng	98

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

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