

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP031622\
 Data File : BP009426.D
 Acq On : 16 Mar 2022 17:12
 Operator : CG/JU
 Sample : N1952-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CON-2

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/17/2022
 Supervised By : Jagrut Upadhyay 03/17/2022

Quant Time: Mar 17 01:09:26 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	7.805	152	482698	20.000	ng	-0.02
21) Naphthalene-d8	10.599	136	1894768	20.000	ng	-0.02
39) Acenaphthene-d10	14.451	164	1070234	20.000	ng	-0.01
64) Phenanthrene-d10	17.210	188	1807123	20.000	ng	-0.01
76) Chrysene-d12	21.339	240	1599099	20.000	ng	0.01
86) Perylene-d12	23.780	264	1838287	20.000	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	3352463	101.379	ng	0.00
7) Phenol-d6	7.005	99	4511584	107.359	ng	0.02
23) Nitrobenzene-d5	8.963	82	2939345	85.443	ng	-0.01
42) 2,4,6-Tribromophenol	15.951	330	1529730	118.372	ng	0.00
45) 2-Fluorobiphenyl	13.075	172	7097640	91.202	ng	-0.01
79) Terphenyl-d14	19.792	244	7793371	91.971	ng	0.00
Target Compounds						Qvalue
60) Diethylphthalate	15.281	149	271094	3.527	ng	98
75) Fluoranthene	19.233	202	339983	2.836	ng	89
78) Pyrene	19.592	202	435293	4.032	ng	# 91
83) Chrysene	21.369	228	237169	2.254	ng	# 67
88) Benzo(b)fluoranthene	23.039	252	500489m	4.032	ng	
90) Benzo(a)pyrene	23.668	252	365980	3.109	ng	# 1
92) Benzo(g,h,i)perylene	27.092	276	241828	2.021	ng	# 46

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

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