

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020822\
 Data File : BP009055.D
 Acq On : 08 Feb 2022 13:17
 Operator : CG/JU
 Sample : N1362-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RO-341

Quant Time: Feb 08 14:05:13 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 07 15:13:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	200690	20.000	ng	0.00
21) Naphthalene-d8	10.604	136	807496	20.000	ng	0.00
39) Acenaphthene-d10	14.445	164	447084	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	791314	20.000	ng	0.00
76) Chrysene-d12	21.304	240	761798	20.000	ng	0.00
86) Perylene-d12	23.698	264	994834	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	958671	85.210	ng	0.00
7) Phenol-d6	6.987	99	1178308	79.483	ng	0.00
23) Nitrobenzene-d5	8.963	82	774044	54.058	ng	0.00
42) 2,4,6-Tribromophenol	15.945	330	465911	78.050	ng	0.00
45) 2-Fluorobiphenyl	13.069	172	1976856	61.899	ng	0.00
79) Terphenyl-d14	19.769	244	1949974	46.032	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	7.016	94	39928	2.679	ng	75
75) Fluoranthene	19.222	202	181336	3.776	ng	97
78) Pyrene	19.575	202	191420	3.532	ng	97
81) Benzo(a)anthracene	21.286	228	111887	2.279	ng	99
83) Chrysene	21.339	228	114904	2.428	ng	97
88) Benzo(b)fluoranthene	22.974	252	198652	3.244	ng	# 93
90) Benzo(a)pyrene	23.592	252	136597	2.675	ng	# 90
92) Benzo(g,h,i)perylene	26.968	276	123423	2.041	ng	# 84

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

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