

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062321\
 Data File : BP006037.D
 Acq On : 23 Jun 2021 16:27
 Operator : CG/JU
 Sample : M2709-10RX
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 0249-AMND-COMP-N(CL-22)RX

Manual Integrations
 APPROVED

mohammad
 6/24/2021 9:51:40 AM

Quant Time: Jun 24 01:16:39 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 22 17:52:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.917	152	69792	20.000	ng	0.00
21) Naphthalene-d8	10.722	136	271474	20.000	ng	# 0.00
39) Acenaphthene-d10	14.546	164	169327	20.000	ng	0.00
64) Phenanthrene-d10	17.292	188	339550	20.000	ng	0.00
76) Chrysene-d12	21.369	240	384621	20.000	ng	0.00
86) Perylene-d12	23.774	264	444123	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	34317	7.890	ng	0.00
7) Phenol-d6	7.093	99	297622	52.793	ng	0.00
23) Nitrobenzene-d5	9.081	82	289051	52.201	ng	0.00
42) 2,4,6-Tribromophenol	16.040	330	11633	4.003	ng	0.00
45) 2-Fluorobiphenyl	13.175	172	698963	54.205	ng	0.00
79) Terphenyl-d14	19.839	244	1192211	58.042	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.334	178	106539	5.225	ng	98
75) Fluoranthene	19.298	202	399915	16.149	ng	99
78) Pyrene	19.645	202	432965	16.125	ng	100
81) Benzo(a)anthracene	21.351	228	350764	12.634	ng	97
83) Chrysene	21.404	228	370817m	13.789	ng	
87) Indeno(1,2,3-cd)pyrene	26.298	276	263416	6.936	ng	98
88) Benzo(b)fluoranthene	23.039	252	537499	17.670	ng	# 97
89) Benzo(k)fluoranthene	23.080	252	176822m	5.995	ng	
90) Benzo(a)pyrene	23.669	252	392195	13.714	ng	# 97
91) Dibenzo(a,h)anthracene	26.298	278	83970	2.569	ng	# 81
92) Benzo(g,h,i)perylene	27.080	276	271520	8.408	ng	98

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

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