

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041221\
 Data File : BP005277.D
 Acq On : 12 Apr 2021 21:23
 Operator : CG/JU
 Sample : M1996-02 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 26878

Manual Integrations
 APPROVED

mohammad
 4/14/2021 11:49:41 AM

Quant Time: Apr 13 00:28:21 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 09 18:44:11 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	42951	20.00	ng	# 0.00
21) Naphthalene-d8	10.469	136	145877	20.00	ng	# 0.00
39) Acenaphthene-d10	14.339	164	93086	20.00	ng	0.00
64) Phenanthrene-d10	17.110	188	206292	20.00	ng	# 0.00
76) Chrysene-d12	21.210	240	203765	20.00	ng	0.00
86) Perylene-d12	23.480	264	208833	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.281	112	70023	17.11	ng	0.00
7) Phenol-d6	6.863	99	75314	16.13	ng	0.00
23) Nitrobenzene-d5	8.840	82	49267	10.65	ng	0.00
42) 2,4,6-Tribromophenol	15.845	330	18935	16.11	ng	0.00
45) 2-Fluorobiphenyl	12.951	172	72476	10.51	ng	0.00
79) Terphenyl-d14	19.686	244	116162	11.20	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.151	178	62917	5.40	ng	# 88
75) Fluoranthene	19.133	202	84094	5.88	ng	94
78) Pyrene	19.486	202	77071	6.09	ng	# 95
81) Benzo(a)anthracene	21.192	228	43419	3.23	ng	93
83) Chrysene	21.245	228	40249m	3.08	ng	
87) Indeno(1,2,3-cd)pyrene	25.804	276	32798	2.05	ng	# 83
88) Benzo(b)fluoranthene	22.792	252	56589m	4.14	ng	
90) Benzo(a)pyrene	23.380	252	42160	3.33	ng	# 94
92) Benzo(g,h,i)perylene	26.521	276	34282	2.55	ng	# 92

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

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