

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041221\
 Data File : BP005266.D
 Acq On : 12 Apr 2021 15:08
 Operator : CG/JU
 Sample : M1967-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SED-1C

Manual Integrations
 APPROVED

mohammad
 4/14/2021 11:49:21 AM

Quant Time: Apr 12 16:09:57 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	26223	20.00	ng	# 0.00
21) Naphthalene-d8	10.469	136	82057	20.00	ng	# 0.00
39) Acenaphthene-d10	14.339	164	48109	20.00	ng	0.00
64) Phenanthrene-d10	17.104	188	114351	20.00	ng	# 0.00
76) Chrysene-d12	21.204	240	149111	20.00	ng	0.00
86) Perylene-d12	23.480	264	159515	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.275	112	288448	115.44	ng	0.00
7) Phenol-d6	6.863	99	304153	106.68	ng	0.00
23) Nitrobenzene-d5	8.840	82	197647	75.96	ng	0.00
42) 2,4,6-Tribromophenol	15.839	330	69748	114.85	ng	0.00
45) 2-Fluorobiphenyl	12.951	172	238304	66.87	ng	0.00
79) Terphenyl-d14	19.680	244	401752	52.93	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.798	163	8794	2.49	ng	# 93
71) Phenanthrene	17.145	178	24539	3.80	ng	98
72) Anthracene	17.239	178	19044	3.04	ng	96
75) Fluoranthene	19.127	202	82537	10.41	ng	99
78) Pyrene	19.480	202	66918	7.22	ng	99
81) Benzo(a)anthracene	21.192	228	52280	5.31	ng	95
83) Chrysene	21.239	228	58485	6.11	ng	98
88) Benzo(b)fluoranthene	22.792	252	65079	6.23	ng	96
89) Benzo(k)fluoranthene	22.827	252	27672m	2.77	ng	
90) Benzo(a)pyrene	23.374	252	47918	4.96	ng	98
92) Benzo(g,h,i)perylene	26.509	276	33370	3.25	ng	96

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

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