

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041321\
 Data File : BP005294.D
 Acq On : 13 Apr 2021 19:08
 Operator : CG/JU
 Sample : M2017-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FISHKILL-SP-1

Manual Integrations
 APPROVED

mohammad
 4/14/2021 11:51:12 AM

Quant Time: Apr 14 02:44:39 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	66911	20.00	ng	# 0.00
21) Naphthalene-d8	10.463	136	231464	20.00	ng	# 0.00
39) Acenaphthene-d10	14.340	164	136723	20.00	ng	0.00
64) Phenanthrene-d10	17.104	188	292102	20.00	ng	# 0.00
76) Chrysene-d12	21.204	240	237959	20.00	ng	0.00
86) Perylene-d12	23.480	264	203937	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.281	112	519849	81.54	ng	0.00
7) Phenol-d6	6.869	99	576243	79.21	ng	0.00
23) Nitrobenzene-d5	8.840	82	390024	53.14	ng	0.00
42) 2,4,6-Tribromophenol	15.845	330	153531	88.96	ng	0.00
45) 2-Fluorobiphenyl	12.951	172	628452	62.05	ng	0.00
79) Terphenyl-d14	19.681	244	881274	72.75	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.798	163	23803	2.37	ng	99
71) Phenanthrene	17.145	178	60706	3.68	ng	97
75) Fluoranthene	19.128	202	154947	7.65	ng	98
78) Pyrene	19.481	202	164586	11.13	ng	98
81) Benzo(a)anthracene	21.192	228	84018	5.35	ng	# 95
83) Chrysene	21.239	228	85178	5.58	ng	# 94
87) Indeno(1,2,3-cd)pyrene	25.815	276	37663	2.42	ng	# 90
88) Benzo(b)fluoranthene	22.792	252	91019	6.82	ng	95
89) Benzo(k)fluoranthene	22.827	252	28892m	2.26	ng	
90) Benzo(a)pyrene	23.380	252	61345	4.97	ng	98
92) Benzo(g,h,i)perylene	26.527	276	37847	2.88	ng	95

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

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