

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP072120\
 Data File : BP002812.D
 Acq On : 21 Jul 2020 14:03
 Operator : CG/JU
 Sample : L3352-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 HS-01

Manual Integrations
 APPROVED

Sohil
 7/24/2020 12:10:19 PM

Quant Time: Jul 21 15:14:28 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 13 10:54:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	130005	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	507321	20.00	ng	-0.01
39) Acenaphthene-d10	14.20	164	309378	20.00	ng	-0.01
64) Phenanthrene-d10	16.96	188	653932	20.00	ng	-0.01
76) Chrysene-d12	21.07	240	546964	20.00	ng	0.00
86) Perylene-d12	23.26	264	626665	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.18	112	1008158	130.84	ng	0.00
7) Phenol-d6	6.75	99	1300592	118.27	ng	0.00
23) Nitrobenzene-d5	8.69	82	813943	85.81	ng	-0.01
42) 2,4,6-Tribromophenol	15.70	330	527206	163.25	ng	0.00
45) 2-Fluorobiphenyl	12.81	172	1714288	85.72	ng	-0.01
79) Terphenyl-d14	19.55	244	2036212	85.35	ng	-0.01
Target Compounds						
50) Dimethylphthalate	13.66	163	244329	10.355	ng	99
71) Phenanthrene	17.00	178	153965	4.349	ng	99
72) Anthracene	17.09	178	73052	2.103	ng	98
75) Fluoranthene	18.99	202	556704	13.570	ng	99
78) Pyrene	19.34	202	565463	14.973	ng	99
81) Benzo(a)anthracene	21.05	228	340112	9.606	ng	99
83) Chrysene	21.10	228	327567	9.719	ng	98
87) Indeno(1,2,3-cd)pyrene	25.46	276	256832	5.670	ng	# 93
88) Benzo(b)fluoranthene	22.60	252	451613	11.606	ng	98
89) Benzo(k)fluoranthene	22.64	252	154517m	4.031	ng	
90) Benzo(a)pyrene	23.16	252	328786	8.916	ng	98
92) Benzo(a,h,i)perylene	26.13	276	255107	6.888	ng	96

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

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