

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100920\
 Data File : BP003596.D
 Acq On : 09 Oct 2020 16:39
 Operator : CG/JU
 Sample : L4333-18
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PM-WC-C

Manual Integrations
 APPROVED

mohammad
 10/12/2020 11:34:51 AM

Quant Time: Oct 09 17:14:16 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 08 15:55:58 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.481	152	43067	20.00	ng	0.00
21) Naphthalene-d8	10.251	136	161319	20.00	ng	# 0.00
39) Acenaphthene-d10	14.145	164	107275	20.00	ng	0.00
64) Phenanthrene-d10	16.910	188	211273	20.00	ng	# 0.01
76) Chrysene-d12	21.010	240	249695	20.00	ng	# 0.00
86) Perylene-d12	23.174	264	281039	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.087	112	32	0.01	ng	-0.01
7) Phenol-d6	6.746	99	5469	1.66	ng	0.06
23) Nitrobenzene-d5	8.634	82	210045	76.49	ng	0.00
42) 2,4,6-Tribromophenol	15.663	330	16	0.01	ng	0.02
45) 2-Fluorobiphenyl	12.751	172	517430	70.54	ng	0.00
79) Terphenyl-d14	19.492	244	819688	68.40	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	10.304	128	89298	10.48	ng	# 95
37) 2-Methylnaphthalene	11.934	142	741242	120.07	ng	# 95
38) 1-Methylnaphthalene	12.157	142	471811	80.49	ng	# 99
46) 1,1'-Biphenyl	12.969	154	134574	16.82	ng	97
49) Acenaphthylene	13.869	152	36702	3.63	ng	# 1
50) Dimethylphthalate	13.610	163	34457	4.02	ng	# 88
52) Acenaphthene	14.210	154	15878m	2.58	ng	
55) Dibenzofuran	14.551	168	112591	11.45	ng	# 22
71) Phenanthrene	16.957	178	849265	73.88	ng	97
75) Fluoranthene	18.945	202	74418	5.15	ng	79
78) Pyrene	19.292	202	185189	11.89	ng	# 95

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

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