

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091820\
 Data File : BP003326.D
 Acq On : 18 Sep 2020 17:41
 Operator : CG/JU
 Sample : L4059-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PAV-B33-091720-0-10

Manual Integrations
 APPROVED

mohammad
 9/21/2020 10:35:50 AM

Quant Time: Sep 18 18:59:08 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 18 14:13:41 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	197069	20.00	ng	0.00
21) Naphthalene-d8	10.375	136	802607	20.00	ng	# 0.00
39) Acenaphthene-d10	14.246	164	484717	20.00	ng	0.00
64) Phenanthrene-d10	17.004	188	1008664	20.00	ng	0.00
76) Chrysene-d12	21.104	240	767452	20.00	ng	0.00
86) Perylene-d12	23.321	264	739576	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.217	112	1241426	110.00	ng	0.00
7) Phenol-d6	6.799	99	1607855	106.88	ng	0.00
23) Nitrobenzene-d5	8.752	82	1055913	77.29	ng	0.00
42) 2,4,6-Tribromophenol	15.745	330	753626	101.94	ng	0.00
45) 2-Fluorobiphenyl	12.869	172	2488029	75.07	ng	0.00
79) Terphenyl-d14	19.586	244	2944543	79.94	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.710	163	532280	13.75	ng	99
71) Phenanthrene	17.045	178	295733	5.39	ng	99
75) Fluoranthene	19.028	202	285943	4.14	ng	99
78) Pyrene	19.381	202	276790	5.78	ng	100
81) Benzo(a)anthracene	21.086	228	113791	2.29	ng	97
83) Chrysene	21.139	228	98493m	2.07	ng	
88) Benzo(b)fluoranthene	22.657	252	118273m	2.53	ng	
90) Benzo(a)pyrene	23.221	252	100031	2.27	ng	# 92

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

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