

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091820\
 Data File : BP003327.D
 Acq On : 18 Sep 2020 18:15
 Operator : CG/JU
 Sample : L4062-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EROC-CF-1

Manual Integrations
 APPROVED

mohammad
 9/21/2020 10:35:46 AM

Quant Time: Sep 19 06:18:48 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	189099	20.00	ng	0.00
21) Naphthalene-d8	10.375	136	774435	20.00	ng	# 0.00
39) Acenaphthene-d10	14.245	164	465733	20.00	ng	0.00
64) Phenanthrene-d10	17.004	188	954627	20.00	ng	0.00
76) Chrysene-d12	21.104	240	722711	20.00	ng	0.00
86) Perylene-d12	23.315	264	694068	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.216	112	1106209	102.15	ng	0.00
7) Phenol-d6	6.799	99	1414567	98.00	ng	0.00
23) Nitrobenzene-d5	8.752	82	926245	70.26	ng	0.00
42) 2,4,6-Tribromophenol	15.751	330	589726	83.02	ng	0.00
45) 2-Fluorobiphenyl	12.869	172	2106143	66.14	ng	0.00
79) Terphenyl-d14	19.586	244	2486004	71.67	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.710	163	211243	5.68	ng	100
75) Fluoranthene	19.027	202	224571	3.44	ng	99
78) Pyrene	19.380	202	221399	4.91	ng	99
81) Benzo(a)anthracene	21.086	228	117760	2.52	ng	98
83) Chrysene	21.139	228	109448	2.44	ng	98
88) Benzo(b)fluoranthene	22.657	252	142011m	3.24	ng	
90) Benzo(a)pyrene	23.221	252	109472	2.65	ng	# 95

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

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