

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091820\
 Data File : BP003328.D
 Acq On : 18 Sep 2020 18:49
 Operator : CG/JU
 Sample : L4062-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EROC-CF-2

Quant Time: Sep 19 06:19:05 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	186955	20.00	ng	0.00
21) Naphthalene-d8	10.381	136	759581	20.00	ng	# 0.00
39) Acenaphthene-d10	14.245	164	462075	20.00	ng	0.00
64) Phenanthrene-d10	17.004	188	936318	20.00	ng	0.00
76) Chrysene-d12	21.104	240	703714	20.00	ng	0.00
86) Perylene-d12	23.321	264	681904	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.216	112	985025	92.00	ng	0.00
7) Phenol-d6	6.799	99	1268785	88.90	ng	0.00
23) Nitrobenzene-d5	8.752	82	834523	64.54	ng	0.00
42) 2,4,6-Tribromophenol	15.751	330	525817	74.61	ng	0.00
45) 2-Fluorobiphenyl	12.869	172	1928035	61.03	ng	0.00
79) Terphenyl-d14	19.586	244	2267649	67.14	ng	0.00
Target Compounds						
						Qvalue
50) Dimethylphthalate	13.710	163	182622	4.95	ng	99
71) Phenanthrene	17.045	178	105686	2.07	ng	99
75) Fluoranthene	19.027	202	224067	3.50	ng	100
78) Pyrene	19.380	202	237409	5.41	ng	99
81) Benzo(a)anthracene	21.086	228	116084	2.55	ng	98
83) Chrysene	21.139	228	108137	2.47	ng	97
88) Benzo(b)fluoranthene	22.657	252	135493	3.15	ng	# 93
90) Benzo(a)pyrene	23.221	252	102648	2.53	ng	96

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

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