

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
 Data File : BP003017.D
 Acq On : 18 Aug 2020 14:53
 Operator : CG/JU
 Sample : L3712-09
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 V-23615

Manual Integrations
 APPROVED

mohammad
 8/19/2020 2:25:01 PM

Quant Time: Aug 18 19:11:55 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 17 14:07:35 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	371684	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	1481696	20.00	ng	0.00
39) Acenaphthene-d10	14.32	164	756864	20.00	ng	0.00
64) Phenanthrene-d10	17.07	188	1722373	20.00	ng	0.00
76) Chrysene-d12	21.17	240	1649761	20.00	ng	0.00
86) Perylene-d12	23.42	264	1715570	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	1505902	62.01	ng	0.00
7) Phenol-d6	6.86	99	1765848	52.37	ng	0.00
23) Nitrobenzene-d5	8.83	82	960073	35.09	ng	0.00
42) 2,4,6-Tribromophenol	15.82	330	621727	66.09	ng	0.00
45) 2-Fluorobiphenyl	12.95	172	2392640	43.19	ng	0.00
79) Terphenyl-d14	19.65	244	3179045	38.58	ng	0.00
Target Compounds						
32) Benzoic acid	9.76	122	3833m	7.920	ng	Qvalue
50) Dimethylphthalate	13.78	163	332271	5.613	ng	99
75) Fluoranthene	19.09	202	264015	2.269	ng	99
78) Pyrene	19.45	202	254478	2.207	ng	97
88) Benzo(b)fluoranthene	22.74	252	254102	2.099	ng	98

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

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