

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG041720\
 Data File : BG045084.D
 Acq On : 19 Apr 2020 2:20
 Operator : CG/JU
 Sample : L2279-07
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OILY-WATER

Manual Integrations
 APPROVED

mohammad
 4/21/2020 8:21:38 AM

Quant Time: Apr 20 05:43:21 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 17 11:08:29 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	63427	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	283671	20.00	ng	0.00
39) Acenaphthene-d10	14.66	164	213892	20.00	ng	0.00
64) Phenanthrene-d10	17.40	188	474217	20.00	ng	0.00
76) Chrysene-d12	21.66	240	421715	20.00	ng	0.00
87) Perylene-d12	24.86	264	452350	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	48221	12.55	ng	0.02
7) Phenol-d6	7.20	99	139604	24.46	ng	0.02
23) Nitrobenzene-d5	9.18	82	230026	37.00	ng	0.00
42) 2,4,6-Tribromophenol	16.15	330	18461m	5.59	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	639090	43.81	ng	0.00
79) Terphenyl-d14	20.01	244	985446	50.93	ng	0.00
Target Compounds						
3) Pyridine	3.89	79	481139	91.524	ng	98
50) Dimethylphthalate	14.11	163	203397	11.679	ng	100
77) Benzidine	19.63	184	118668	4.333	ng	96

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

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