

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081720\
 Data File : BP003002.D
 Acq On : 17 Aug 2020 14:54
 Operator : CG/JU
 Sample : L3688-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BRICK-PILE

Manual Integrations
 APPROVED

mohammad
 8/19/2020 2:20:37 PM

Quant Time: Aug 18 01:52:28 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	300338	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	1363905	20.00	ng	0.00
39) Acenaphthene-d10	14.32	164	757440	20.00	ng	0.00
64) Phenanthrene-d10	17.07	188	1653511	20.00	ng	0.00
76) Chrysene-d12	21.17	240	1929714	20.00	ng	0.00
86) Perylene-d12	23.42	264	1921633	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	1131069	57.64	ng	0.00
7) Phenol-d6	6.86	99	1458078	53.52	ng	0.00
23) Nitrobenzene-d5	8.83	82	777945	30.89	ng	0.00
42) 2,4,6-Tribromophenol	15.82	330	691230	73.42	ng	0.00
45) 2-Fluorobiphenyl	12.95	172	2256628	40.71	ng	0.00
79) Terphenyl-d14	19.65	244	3313382	34.37	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.78	163	329139	5.556	ng	99
71) Phenanthrene	17.12	178	391362	4.036	ng	99
75) Fluoranthene	19.09	202	1152155	10.313	ng	99
78) Pyrene	19.45	202	1018534	7.550	ng	98
81) Benzo(a)anthracene	21.15	228	668910	5.089	ng	100
83) Chrysene	21.20	228	640433	5.165	ng	98
87) Indeno(1,2,3-cd)pyrene	25.69	276	403982	2.625	ng	96
88) Benzo(b)fluoranthene	22.74	252	866650m	6.391	ng	
89) Benzo(k)fluoranthene	22.77	252	256460m	2.001	ng	
90) Benzo(a)pyrene	23.32	252	562250	4.589	ng	100
92) Benzo(g,h,i)perylene	26.39	276	378610	2.989	ng	98

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

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