

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP061720\
 Data File : BP002433.D
 Acq On : 17 Jun 2020 23:06
 Operator : CG/JU
 Sample : L3024-19
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RO-17

Manual Integrations
 APPROVED

mohammad
 6/18/2020 11:54:57 AM

Quant Time: Jun 18 01:38:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 05 15:08:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	147156	20.00	ng	0.00
21) Naphthalene-d8	10.36	136	620400	20.00	ng	0.00
39) Acenaphthene-d10	14.23	164	383661	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	853099	20.00	ng	-0.01
76) Chrysene-d12	21.09	240	870262	20.00	ng	-0.02
86) Perylene-d12	23.29	264	905667	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	884452	111.22	ng	0.00
7) Phenol-d6	6.78	99	1223689	115.57	ng	0.00
23) Nitrobenzene-d5	8.73	82	822197	79.14	ng	-0.01
42) 2,4,6-Tribromophenol	15.73	330	427771	116.45	ng	0.00
45) 2-Fluorobiphenyl	12.85	172	1768436	77.50	ng	0.00
79) Terphenyl-d14	19.58	244	2740243	78.96	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.70	163	217436	8.271	ng	99
78) Pyrene	19.36	202	129478	2.429	ng	99
83) Chrysene	21.13	228	104900	2.229	ng	94
88) Benzo(b)fluoranthene	22.63	252	280749m	5.748	ng	
90) Benzo(a)pyrene	23.19	252	166022	3.627	ng	96

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

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