

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062420\
 Data File : BP002527.D
 Acq On : 24 Jun 2020 17:35
 Operator : CG/JU
 Sample : L3089-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RB47185-RT

Manual Integrations
 APPROVED

mohammad
 6/25/2020 1:07:32 PM

Quant Time: Jun 24 18:28:00 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 24 13:54:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	237871	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	953993	20.00	ng	0.00
39) Acenaphthene-d10	14.23	164	547219	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	1071580	20.00	ng	0.00
76) Chrysene-d12	21.09	240	796112	20.00	ng	0.00
86) Perylene-d12	23.29	264	904134	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	1439677	111.99	ng	0.00
7) Phenol-d6	6.78	99	1916801	112.00	ng	0.00
23) Nitrobenzene-d5	8.73	82	1332929	83.44	ng	0.00
42) 2,4,6-Tribromophenol	15.73	330	635026	121.21	ng	0.00
45) 2-Fluorobiphenyl	12.85	172	2653401	81.53	ng	0.00
79) Terphenyl-d14	19.57	244	3019009	103.52	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.69	163	359862	9.597	ng	99
71) Phenanthrene	17.03	178	249208	4.972	ng	99
75) Fluoranthene	19.01	202	596239	9.965	ng	98
78) Pyrene	19.36	202	510187	10.462	ng	98
81) Benzo(a)anthracene	21.07	228	244228	5.374	ng	98
83) Chrysene	21.13	228	278099m	6.459	ng	
87) Indeno(1,2,3-cd)pyrene	25.49	276	180995	3.201	ng	99
88) Benzo(b)fluoranthene	22.63	252	369095	7.570	ng	# 97
89) Benzo(k)fluoranthene	22.67	252	133719m	2.886	ng	
90) Benzo(a)pyrene	23.19	252	245006	5.362	ng	98
92) Benzo(g,h,i)perylene	26.16	276	171741	3.655	ng	96

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

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