

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062420\
 Data File : BP002528.D
 Acq On : 24 Jun 2020 18:09
 Operator : CG/JU
 Sample : L3089-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RB44767-RT

Manual Integrations
 APPROVED

mohammad
 6/25/2020 1:07:35 PM

Quant Time: Jun 24 18:42:01 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	212618	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	806044	20.00	ng	0.00
39) Acenaphthene-d10	14.23	164	450877	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	878431	20.00	ng	0.00
76) Chrysene-d12	21.09	240	794888	20.00	ng	0.00
86) Perylene-d12	23.29	264	910627	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	1456246	126.74	ng	0.00
7) Phenol-d6	6.78	99	1903888	124.45	ng	0.00
23) Nitrobenzene-d5	8.73	82	1309283	97.00	ng	0.00
42) 2,4,6-Tribromophenol	15.73	330	593921	137.58	ng	0.00
45) 2-Fluorobiphenyl	12.85	172	2527735	94.26	ng	0.00
79) Terphenyl-d14	19.57	244	2878841	97.01	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.69	163	243485	7.881	ng	100
71) Phenanthrene	17.03	178	195745	4.764	ng	99
75) Fluoranthene	19.01	202	525399	10.712	ng	98
78) Pyrene	19.36	202	451224	9.267	ng	99
81) Benzo(a)anthracene	21.07	228	259515	5.720	ng	97
83) Chrysene	21.13	228	309569m	7.201	ng	
87) Indeno(1,2,3-cd)pyrene	25.49	276	205455	3.607	ng	99
88) Benzo(b)fluoranthene	22.63	252	426924	8.694	ng	99
89) Benzo(k)fluoranthene	22.67	252	159192m	3.412	ng	
90) Benzo(a)pyrene	23.19	252	274841	5.972	ng	96
92) Benzo(g,h,i)perylene	26.16	276	195745	4.137	ng	99

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

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