

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061621\
 Data File : BP005951.D
 Acq On : 16 Jun 2021 15:20
 Operator : CG/JU
 Sample : M2678-07RX
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 0241-AMND-COMP-F(CL-05)RX

Manual Integrations
 APPROVED

mohammad
 6/17/2021 12:16:26 PM

Quant Time: Jun 16 16:05:51 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 16 11:35:30 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	75796	20.00	ng	0.00
21) Naphthalene-d8	10.734	136	291539	20.00	ng	0.00
39) Acenaphthene-d10	14.557	164	186323	20.00	ng	0.00
64) Phenanthrene-d10	17.304	188	391452	20.00	ng	0.00
76) Chrysene-d12	21.374	240	459569	20.00	ng	0.00
86) Perylene-d12	23.786	264	529984	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.505	112	4144	0.88	ng	0.00
7) Phenol-d6	7.105	99	88710	14.49	ng	0.00
23) Nitrobenzene-d5	9.099	82	356580	59.96	ng	0.00
42) 2,4,6-Tribromophenol	16.057	330	680	0.21	ng	0.00
45) 2-Fluorobiphenyl	13.187	172	807290	56.89	ng	0.00
79) Terphenyl-d14	19.851	244	1407812	57.36	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	17.345	178	230867	9.82	ng	99
72) Anthracene	17.439	178	138245	5.98	ng	99
75) Fluoranthene	19.304	202	721514	25.27	ng	98
78) Pyrene	19.657	202	639233	19.92	ng	99
81) Benzo(a)anthracene	21.357	228	455342	13.73	ng	98
83) Chrysene	21.410	228	397958m	12.39	ng	
87) Indeno(1,2,3-cd)pyrene	26.315	276	223573	4.93	ng	98
88) Benzo(b)fluoranthene	23.051	252	465065	12.81	ng	97
89) Benzo(k)fluoranthene	23.092	252	177959m	5.06	ng	
90) Benzo(a)pyrene	23.680	252	378898	11.10	ng	99
92) Benzo(g,h,i)perylene	27.098	276	195361	5.07	ng	98

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

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