

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062921\
 Data File : BP006104.D
 Acq On : 29 Jun 2021 22:36
 Operator : CG/JU
 Sample : M2884-01 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OK-03-062821

Manual Integrations
 APPROVED

mohammad
 6/30/2021 2:14:07 PM

Quant Time: Jun 30 01:14:48 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 28 11:34:30 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	144321	20.000	ng	0.00
21) Naphthalene-d8	10.704	136	533834	20.000	ng	0.00
39) Acenaphthene-d10	14.534	164	313745	20.000	ng	0.00
64) Phenanthrene-d10	17.281	188	607571	20.000	ng	0.00
76) Chrysene-d12	21.351	240	527960	20.000	ng	0.00
86) Perylene-d12	23.757	264	551439	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	332418	36.960	ng	0.00
7) Phenol-d6	7.087	99	400743	34.376	ng	0.00
23) Nitrobenzene-d5	9.069	82	244193	22.427	ng	0.00
42) 2,4,6-Tribromophenol	16.028	330	182329	33.864	ng	0.00
45) 2-Fluorobiphenyl	13.157	172	570591	23.881	ng	0.00
79) Terphenyl-d14	19.828	244	797011	28.267	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.322	178	143221	3.925	ng	99
75) Fluoranthene	19.281	202	361425	8.157	ng	97
78) Pyrene	19.633	202	276887	7.512	ng	99
81) Benzo(a)anthracene	21.339	228	131496	3.450	ng	97
83) Chrysene	21.386	228	149747m	4.057	ng	
87) Indeno(1,2,3-cd)pyrene	26.280	276	99208	2.104	ng	99
88) Benzo(b)fluoranthene	23.022	252	196246	5.196	ng	# 90
90) Benzo(a)pyrene	23.645	252	124101	3.495	ng	# 90
92) Benzo(g,h,i)perylene	27.062	276	103243	2.575	ng	# 95

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

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