

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071221\
 Data File : BP006189.D
 Acq On : 12 Jul 2021 19:39
 Operator : CG/JU
 Sample : M3006-01 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CL-02-070921

Manual Integrations
 APPROVED

mohammad
 7/13/2021 1:42:21 PM

Quant Time: Jul 13 04:16:31 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 12 14:59:51 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.864	152	239718	20.000	ng	0.00
21) Naphthalene-d8	10.663	136	967329	20.000	ng	0.00
39) Acenaphthene-d10	14.498	164	552067	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	1015403	20.000	ng	# 0.00
76) Chrysene-d12	21.328	240	912298	20.000	ng	# 0.00
86) Perylene-d12	23.710	264	1029741	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	817773	58.698	ng	0.00
7) Phenol-d6	7.052	99	1022912	55.508	ng	0.00
23) Nitrobenzene-d5	9.028	82	571719	37.332	ng	0.00
42) 2,4,6-Tribromophenol	15.993	330	389446	59.252	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	1404317	37.605	ng	0.00
79) Terphenyl-d14	19.798	244	1742905	37.667	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	17.292	178	207443	3.617	ng	97
75) Fluoranthene	19.257	202	327923	5.217	ng	96
78) Pyrene	19.604	202	308431	5.002	ng	98
81) Benzo(a)anthracene	21.310	228	170992	2.963	ng	97
83) Chrysene	21.363	228	148015m	2.596	ng	
88) Benzo(b)fluoranthene	22.980	252	226149	3.620	ng	# 46
90) Benzo(a)pyrene	23.604	252	171997	2.934	ng	# 47

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

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