

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071321\
 Data File : BP006212.D
 Acq On : 13 Jul 2021 21:41
 Operator : CG/JU
 Sample : M2997-01DL 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RBR-200029DL

Manual Integrations
 APPROVED

mohammad
 7/15/2021 11:30:35 AM

Quant Time: Jul 14 01:54:05 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	7.852	152	189953	20.000	ng	-0.01
21) Naphthalene-d8	10.651	136	786332	20.000	ng	#-0.01
39) Acenaphthene-d10	14.487	164	427611	20.000	ng	-0.01
64) Phenanthrene-d10	17.239	188	815604	20.000	ng	#-0.01
76) Chrysene-d12	21.333	240	889022	20.000	ng	# 0.00
86) Perylene-d12	23.733	264	1065749	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	115344	10.448	ng	0.00
7) Phenol-d6	7.057	99	139116	9.527	ng	0.01
23) Nitrobenzene-d5	9.034	82	84771	6.810	ng	0.00
42) 2,4,6-Tribromophenol	15.986	330	56791	11.155	ng	-0.01
45) 2-Fluorobiphenyl	13.116	172	223356	7.722	ng	0.00
79) Terphenyl-d14	19.798	244	342174	7.589	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	17.280	178	95349	2.070	ng	# 93
75) Fluoranthene	19.251	202	165175	3.272	ng	85
78) Pyrene	19.604	202	149882	2.494	ng	# 83
84) Bis(2-ethylhexyl)phtha...	21.227	149	1630512	37.063	ng	93
88) Benzo(b)fluoranthene	23.010	252	177707m	2.749	ng	

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

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